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Transient Thermal Modeling of the Nonscanning ERBE Detector

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Transient Thermal Modeling of the
Nonscanning ERBE Detector

1. Introduction

This report summarizes the activity at VPI under Contract NAS1-16508 between 13 November 1980 and 30 September 1982. During this period a numerical model of the ERBE wide-field-of-view total radiometer channel was developed, and the model was applied to the TRW calibration procedure. Other activities included consulting with NASA and contractor scientists and engineers on the design and calibration of radiometric instruments and on the interpretation of data from such instruments.

Three major documents have resulted from this work: a mid-term report entitled, Application of the Monte Carlo Method to the Transient Thermal Modeling of a Diffuse-Specular Radiometer Cavity (Leo Eskin's M.S. thesis)¹, a Users' Manual for the cavity transient thermal model program (included as an appendix to reference 1, and the TRW Calibration Simulation Users' Manual, the final revised version which appears in the Appendix of this final report. In addition, a paper was presented at the Fourth AMS Conference on Atmospheric Radiation in Toronto in June of 1981².

The final numerical model of the ERBE instrument, the "TRW Calibration Simulation Model", consists of two sub-models which interact with each other within the structure of the overall model. One of these sub-models computes the transient thermal response of the cone

cavity, and the other computes the transient thermal response of the instrument body itself. The details of the cavity sub-model are described in reference 1 and thus are not reiterated here.

In the present report we document the instrument body sub-model and its interaction with the cavity sub-model and its surroundings. We indicate in some detail how this model can be modified to take into account the "two-temperature" aspect of the TRW calibration source with its mounting collar, or baffle. We also outline how the "TRW Calibration Model" can be modified and extended to serve as a flight simulation model, including operation as a visible channel. Finally, we suggest a procedure for using the model to establish a real-time interpretation strategy.

2. The TRW Calibration Simulation Model

The Users' Manual for the "TRW Calibration Simulation Model" appears in the Appendix of this report. This is the final revised version originally transmitted to NASA Langley by Dave Parekh in September, 1982. At that time Mr. Parekh also sent NASA a magnetic tape containing this model. In this section we describe the parts of this model not already documented in Reference 1. This is also where we suggest the modifications referred to in the Introduction.

2.1. TRW FRONT

"TRW FRONT" is a subprogram of the TRW Calibration Simulation Model in which Monte Carlo techniques are used to compute the

distribution factors for radiation among the surfaces of the "front end" of the instrument, between these surfaces and the blackbody calibration source, and between these surfaces and the cavity. By "front end", we mean the field-of-view (FOV) limiter and the surface in the plane of the precision aperture to which the FOV limiter is attached. The general principles of the Monte Carlo technique are described in great detail in Reference 1. Because the logic of "TRW FRONT" is very similar to that of the cavity distribution factor program documented in Reference 1, we will not reiterate the details here. However, the program listing, with notes, appears on pages 47-76 of the Appendix.

The case documented in the Appendix is for the wide field-of-view total channel. It can be easily modified for any other geometry by changing values of the data read by the program, "READ(5,97)", page 49. These data appear after the subroutines which serve "TRW FRONT", on page 75 of the Appendix. Each element of this data set is carefully labeled on page 76, described on page 47, and referred to in Fig. 1 of the February, 1982, Progress Report.

2.1.1. Modification to Account for a Two-Temperature Calibration Source

The present version of "TRW FRONT" assumes that the instrument faces a uniform temperature blackbody source. The blackbody calibration source is thus modeled as a perfectly absorbing sector in the plane of the FOV limiter. This is exactly equivalent to any real

geometry black cavity into which the instrument might face as long as the cavity is isothermal. If, however, as in the case of the actual TRW calibration source, the calibration cavity is composed of two surfaces having different temperatures, "TRW FRONT" must be modified to take into account its actual shape. The page-by-page procedure for this modification is outlined below:

1. Page 49 : change $NN3 = NN2 + 3$

to $NN3 = NN2 + 4$

change $MM3 = M3 + 1$

to $MM3 = M3 - NTH$

2. Page 50 : change

GOTO(101, 102,..., 110),J

101 CONTINUE

...

102 CONTINUE

to

GOTO(100, 101, 102,..., 110),J

100 CONTINUE

statements similar to those presently between "101 CONTINUE" and "102 CONTINUE" which establish the x, y and z coordinants of a randomly selected source location on the controlled surface of the calibration cavity

101 CONTINUE

statements which establish the x, y and z coordinants of a randomly selected source location on the concentric collar, or baffle, which couples the calibration source to the ERBE instrument

102 CONTINUE

3. Page 55: change IF(TZ1.EQ.H2) GOTO 201

to IF(TZ1.GT.H2) GOTO 201

4. Page 56 : change three statements between "201 CONTINUE" and "GOTO 23" to the logic necessary to compute values of UNX, UNY and UNZ, depending on whether the source point is on the blackbody surface itself or on the baffle. This logic depends on the geometry of the calibration source and its interface with the instrument, but should be easy to implement.

5. Page 61 : After "11 CONTINUE" the logic must be changed depending on the geometry of the calibration source/radiometer interface. First, determine if the source point is on the calibration blackbody surface or on the baffle by testing the value of TZI in a manner similar to that on page 60. Then compute TXJ and TYJ in a manner similar to that in the present version of the program. If the source point is on the blackbody surface, begin the search to see if the energy packet has entered the instrument field of view by comparing the value of

$$DSQRT(TXJ*TXJ + TYJ*TYJ)$$
 with the radius of the FOV limiter at TZJ = H2. If this value exceeds the radius of the FOV limiter, the ray is reabsorbed somewhere within the calibration source, in which case go to "RETURN". However, if this value is

less than the radius of the FOV limiter at $TZJ = H2$, the energy packet has entered the instrument field of view, in which case the search continues as presently organized under "11 CONTINUE". The logic is similar if the source point is on the baffle rather than the blackbody surface itself.

2.1.2. Adaptation to a "Real" Source Field

If the "TRW Calibration Simulation Model" is to be modified as a flight simulator, the calibration blackbody source must be replaced with a "real" source field. This real field must, of course, be time varying. It is impractical to work in terms of distribution factors from the earth scene to the surfaces of the instrument, since these factors would all be essentially zero (the fraction of radiation emitted diffusely from Kansas which is absorbed by a 5-mm square surface in earth orbit is too small to compute). We also have demonstrated that an artificial "grid" placed over the FOV inlet plane to represent the earth scene is unsatisfactory because of its "focusing" effect. The most efficient way to adapt the existing model to a "real" field, such as GENDAT, is outlined below:

1. Use GENDAT (or some equivalent source field model) to compute the flux incident to the disk defined by the FOV limiter, as well as to the disk defined by the precision aperture. These calculations are both within the present capabilities of GENDAT.

2. The difference between these two values is the flux incident to the FOV limiter and the "floor" surrounding the precision aperture. Assume that none of this energy is specularly reflected into the cone cavity, in keeping with the design principle of the FOV limiter. The fraction of this energy which is diffusely reflected into the precision aperture is

$$\sum_{i=1}^n D_{i-A} \rho_i^d ,$$

where D_{i-A} is the distribution factor from the i th front end surface element to the disk defined by the precision aperture, computed in "TRW FRONT", and ρ_i^d is the diffuse component of the reflectivity of this element. Note that this calculation assumes that the energy from the scene incident to the front end optics is uniformly distributed across the FOV limiter and floor. This assumption is of minor consequence in view of the fact that it is approximately true for any realistic earth scene, and that the result in any case is a small contribution to the total energy input to the cavity.

3. The fraction of the flux entering the cavity directly from the earth scene which is absorbed by cavity surface element j is

$$D_{A-j} = \epsilon_j \frac{A_j}{A_A} D_{j-A} ,$$

where A_j is the surface area of cavity element j , ϵ_j is its diffuse emissivity, A_A is the precision aperture opening area, and D_{j-A} is the distribution factor from cavity element j to the precision aperture opening computed in "TRW NODE". This calculation assumes that the flux from the earth scene incident to the precision aperture opening is diffuse. Although this assumption is not strictly valid, the error associated with it will be negligible in view of the cavity properties.

2.1.3. Inclusion of a Filter Dome

The extension of the "TRW Calibration Simulation Model" to include the visible channel would be a major modification. It would involve altering the paths of the energy packets which intercepted the filter dome as follows:

1. Upon the intersection of a ray with the filter dome, a random number would be generated to determine if it was reflected. If not reflected, another random number would be generated to see if it was absorbed or transmitted.
2. If the energy packet is absorbed, the dome element would be treated like any other surface element; that is, the counter for the distribution factor from the source element to the dome element would be incremented.
3. If the energy packet is transmitted, diffraction can be included according to Snell's law.
4. If the energy packet is reflected, we have to generate

another random number to determine if the reflection is specular or diffuse.

- a) If specular, treat the dome element like any other surface.
- b) If diffuse, the possibility of forward scattering should be included; that is, uniformly distributed scattering in 4π space rather than 2π space.

After reflection, or scattering, the energy packet is traced to its next surface just as in the case without a dome.

2.2. TRW TEMP

This subprogram computes the transient temperature distributions in the active cavity and body of the instrument based on the distribution factors and conductances computed in the previous subprograms (or user generated in the case of the body elements). As presently configured, this calculation is based on an assumed isothermal black-body source. Arbitrary heat sink and mounting beam temperatures can be introduced to correspond to actual calibration conditions. A detailed annotated listing of this subprogram appears in the Appendix, pages 77-98. The data listings appear on pages 99-104.

The thermal mass of a typical instrument body node is much larger than that of a typical cavity node. For this reason the temperatures of the body nodes change much more slowly than those of the cavity nodes. Thus, we can realize a certain efficiency by using a "two-timing" technique to compute the transient temperature distributions in the instrument body and on the active cavity surfaces. The body

node temperature calculations are performed only after every tenth cavity node temperature calculation. The radiation heat transfer from the FOV limiter to the active cavity used in each cavity temperature calculation is then based on the quasi-steady temperatures of the FOV limiter nodes. The conduction heat transfer from the cavity nodes to the body nodes is also calculated based on the quasi-steady body node temperatures and stored in a buffer until needed for the next body node temperature calculation, at which time this accumulated energy is distributed to the appropriate body nodes.

2.2.1. Detailed Description of the Body Node Temperature Calculation

The active cavity node temperature calculation is well documented in Reference 1. In this section the body node temperature calculation and its interaction with the cavity node temperature calculation is described.

The values of certain constants, defined on page 77, are established on page 80. In particular, the user must change "HST", "TCLAMP", "HEAT" and "TSOURC" to conform to the actual calibration conditions. The statements on pages 80-86 establish the values of the constants, coefficients, conductances, distribution factors, and initial temperatures for the calculations. The actual calculations begin on page 86 with the "280" DO-loop, which extends to the end of the subprogram.

The first step in this sequence is to check to see if it is time to do a body temperature calculation ("IF(KOUNT.EQ.10) GOTO 678", page 87). The actual body temperature calculations lie between

"678 CONTINUE", page 87 , and "679 CONTINUE", page 88 . Between "219 CONTINUE", page 87 , and "GOTO 219", page 88 , an estimate of the new node temperatures is computed based on the old node temperatures, the input from the cavity, the input from the field (calibration source), and the inputs from the heat sink and mounting beam. This calculation loop is repeated until convergence is obtained to a cavity temperature distribution whose rms value no longer changes from one loop to the next. In this loop, "TGBODY" represents the "guess" for the temperature distribution, and "TUBODY" represents the updated temperature distribution. At the end of each cycle through this loop, the guess is updated and, if there is a significant difference between the update and the preceding guess, the loop is repeated. In the present version of the subprogram a change of one part in a million is tolerated in the rms difference between the last two iterations. This tolerance can be changed by the user in the statement "IF (TEST.LE.0.000001) GOTO 241", page 88.

The updated temperature distribution is actually computed in the "230" DO-loop, pages 87-88. The logic in this loop is "hard wired" in that, if the body shape or node configuration is changed, the logic in this loop must be changed accordingly. These changes would be made in the statements between "220 CONTINUE" and "637 CONTINUE", page 87. Between "637 CONTINUE", page 87, and "230 CONTINUE", page 88, the actual temperature calculation is made based on Newton's method. From the statement "230 CONTINUE", page 88, down to the statement "IF (TEST.LE.0.000001) GOTO 241", page 88, the relative rms change in the temperature distribution is calculated and tested.

If the temperature distribution has converged, the body node temperatures "T1BODY" are changed to their new values, "TUBODY", in the "260" DO-loop, page 88. The clock is updated and the new body temperature distributions are written out in the statements which immediately follow. Finally, the power input to the cavity by radiation from the FOV limiter, "QBOD", is computed between "246 CONTINUE" and "679 CONTINUE", page 88.

The statements after "679 CONTINUE" in which the cavity node temperatures are computed, are for the most part identical to those documented in Reference 1. The main differences are on page 91, where the heat input from the front end, "QBOD", is added ("IF (I.LE.4) Q2 = Q2 + QBOD"), and on page 93, where the heat conduction "QTOT" is buffered from the cavity to the body nodes ("IF(IJK.LE.4) QTOT(LL) = QTOT(LL) + ...").

2.2.2. Modifications for the Addition of a Filter Dome

If a filter dome is installed, additional body nodes must be provided. The conductances and distribution factors associated with these new nodes must be calculated and included in the data files. The calculation of the distribution factors would be carried out in "TRW FRONT" as discussed in Section 2.1.3. However, the conductances must be provided by the user. What follows is a page-by-page outline of the modifications required for the inclusion of the filter dome.

1. Page 80: Increase all dimensions "104" to a number which includes the filter dome nodes, and change all dimensions "106" to this new number plus 2 (the number "106" presently refers to the cavity aperture).

Increase the dimension of "SOURCE" to $41 + n$, where n is the number of filter dome nodes.

Increase the value of "NBODY" to include the additional filter dome nodes.

2. Page 82: Change the upper limit of the "601" DO-loop to include the filter dome nodes (that is, increase 41 to $41 + n$, where n is the total number of filter dome nodes).

3. Page 83: Change the "105" in "IF(J.LT.5) JJ = 105" to the total number of body plus filter dome nodes + 1.

Change "IF(J.EQ.41) JJ = 106" to "IF(J.EQ.number1) JJ = number2", where $\text{number1} = 41 + n$ (n being the number of filter dome nodes) and $\text{number2} = \text{total number of body plus filter dome nodes} + 2$.

After "IF(J.GE.33.AND.J.LT.41) etc.", add a similar statement which converts the dome node position indices (established in the data set) to node numbers.

Change the upper limit in the two "401" DO-loops to the total number of body plus filter dome nodes.

4. Page 87: Between "220 CONTINUE" and "636 CONTINUE", add statements similar to those for the existing body nodes, for the filter dome nodes.

5. Page 88: Change the upper limit of the "246" and "661" DO-loops to the total number of body plus filter dome nodes.

In the line above "661 CONTINUE", change "106" to the total number of body plus filter dome nodes + 2.

6. Pages 100,101 & 104: Add thermal physical properties and conductance data (user generated) for filter dome nodes.

3. Use of the Model to Establish a Real-Time Flight Data Interpretation Strategy

The "TRW Calibration Simulation Model" could be used to establish a real-time flight data interpretation strategy. The first step would be to adjust the model (with the filter dome modifications for the visible channel) until it is in good agreement with the actual instrument in the calibration environment. The next step would be to implement the modifications outlined above to convert the model into a flight simulator. Finally, the "calibrated" model would be used to establish an empirical relation between the actual source energy flux at the instrument, computed independently of the instrument response using, for example, GENDAT, the active cavity heater power signal, and one or more (the more the better) key instrument temperatures. It would be hoped that these "key instrument temperatures" would play only secondary roles in this relation. A logical and yet relatively simple form, consistent with Figs. 1 and 2 of the July, 1982, Progress Report, is

suggested by the superposition theorem³ :

$$Q_{\text{heater}}(t) = A \int_0^t \frac{\partial}{\partial t} Q_{\text{source}}(t-\tau) d\tau + B \int_0^t \frac{\partial}{\partial t} T_{\text{HS}}(t-\tau) d\tau \\ + C \int_0^t \frac{\partial}{\partial t} T_c(t-\tau) d\tau .$$

In this relation, $Q_{\text{heater}}(t)$ is the instantaneous heater power (or at least the "count" which represents this power), $Q_{\text{source}}(t-\tau)$ is the flux incident to the precision aperture from the source a time delay τ earlier, $T_{\text{HS}}(t-\tau)$ is the heat sink temperature a time delay τ earlier, and $T_c(t-\tau)$ is the temperature of the mounting beam a time delay τ earlier. The sensitivity coefficients A, B, and C would be obtained by exercising the calibrated model. This relation is inconvenient because Q_{source} appears under an integral. However, it should be possible to write an algorithm capable of inverting this integral in real time.

It is possible that a more complicated model involving more key temperatures will be required. In any case, the numerical model, modified and calibrated as described in this report, should prove very useful in defining this relation.

Once flight data are obtained and interpreted by some data interpretation strategy, the result can be introduced into the complete model to verify that the correct count is indeed obtained.

4. References

1. Eskin, L.D., Application of the Monte Carlo Method to the Transient Thermal Modeling of a Diffuse-Specular Radiometer Cavity, Master's Thesis, Virginia Polytechnic Institute and State University, Blacksburg, Virginia, August 1981.
2. Mahan, J. R., and L. D. Eskin, "Application of Monte Carlo Techniques to Transient Thermal Modeling of Cavity Radiometers Having Diffuse-Specular Surfaces," Fourth American Meteorological Society Conference on Atmospheric Radiation, Toronto, Ontario, Canada, June 16-18, 1981.
3. Eckert, E. R. G., and R. M. Drake, Jr., Analysis of Heat and Mass Transfer (McGraw-Hill, New York, 1972), p. 170.

5. APPENDIX

Transient Thermal Modeling of the
Nonscanning ERBE Detector

TRW Calibration Simulation

USERS MANUAL

Final Revision: March 1983

1. TRW SHAPE

This program subdivides a TRW-type cavity into $NN2 \times NTH$ elements, where $NN2$ is the number of divisions along the cavity axis and NTH is the number of circumferential divisions. When the cavity has been subdivided into elements, the node surface areas and thermal capacities and the inter-node thermal conductances are computed and stored in online disk files.

This program requires two previously generated data file areas on system disk. The first file name should be named A51961.DATA10 and be of sufficient size to hold $M2 + 1$ card images. This file will be written to as unit 10. The second file should be named A51961.DATA15 and should be of sufficient size to hold $M2 \times M2 + NN2 + 2$ card images. This file will be written to as unit 15.

1.1 Nomenclature

ZTB	Matrix of the z-coordinate for each node level top boundary. Must be dimensioned $NN2$.
DZ	Matrix of the z-coordinate width of each node level. Must be dimensioned $NN2$.
AA	Matrix of the node surface areas (m^2). Must be dimensioned $M2$.
CC	Matrix of the node thermal capacities (W-s/K). Must be dimensioned $M2$.
VX(I,I)	X-position of the Ith node in Cartesian coordinates.
VY(I,I)	Y-position of the Ith node in Cartesian coordinates.
VZ(I,I)	Z-position of the Ith node in Cartesian coordinates.
VX(I,J)	"Half-conductance" between nodes I and J where $I > J$.
VY(J,I)	"Half-conductance" between nodes I and J where $J > I$.

(Note: VX, VY, and VZ must all be dimensioned $M2$ by $M2$.)

A	Distance from plane of precision aperture to junction of barrel and cone (mm).
B	Distance from plane of precision aperture to apex of cone (mm).
C	Radius of barrel (mm).

(Note: For the purpose of stating these dimensions, it is assumed that there is no "space ring" gap between the precision aperture and the cavity opening.)

ALPHA	Slope of cone (-).
PI	Ratio of circumference to diameter of a circle (-).
COND	Thermal conductivity (W/m-K).
DENS	Mass density (kg/m**3).
SPEC	Specific heat (W-s/kg-K).
THCK	Thickness (mm).
NARSZ	Number of columns (or rows) in arrays VX, VY, and VZ. This must be the dimensioned value.
NTH	The number of circumferential divisions desired.
FNTH	Floating point version of "NTH".
N1	Number of axial divisions of barrel.
N2	Number of axial divisions of cavity.
FN1, FN2	Floating point versions of N1 and N2.
NN1	Total number of axial divisions through barrel.
NN2	Total number of axial divisions of cavity.
M1	Total number of cylinder nodes.
M2	Total number of cavity nodes.
I	Node number.
J	Index of node level.
FJ	Real value of J.
TZ	Temporary value of z-coordinate of node level J.
TZ1	Temporary value of z-coordinate of node level J-1.
TZ2	Temporary value of z-coordinate of node level J+1.
GK1	"Half-conductance" to node to the "left" of a node on level J.
GK2	"Half-conductance" to node "above" a node on level J, i.e., on level J-1.
GK3	"Half-conductance" to node to the "right" of a node on level J; note that GK3 = GK1.
GK4	"Half-conductance" to node "below" a node on level J, i.e., on level J+1.
TZB1	Temporary z-coordinate of the boundary between nodes on level J and nodes on level J-1.
TZB2	Temporary z-coordinate of the boundary between nodes on level J and nodes on level J+1.
R	Radial location of nodes at node level J.
RB1	Radial position of boundary between nodes on level J and nodes on level J-1.
RB2	Radial position of boundary between nodes on level J and nodes on level J+1.
AREA	Area occupied by a node at node level J (m**2).
J	Auxillary counter defined in "10" DO-loop used to determine which segment of the cavity is being considered.
FJ	Floating point version of J.
TZ	Temporary z-coordinate of node; later converted to a permanent value, VZ(I,I).
R	Radial coordinate of node; later resolved into x- and y-coordinates, VX(I,I) and VY(I,I).
THETA	Circumferential angular location of node.

1.2 Description of Program

It is necessary to specify double precision when this program is compiled on IBM equipment. This step may not be necessary with other processors.

```
IMPLICIT REAL*8 (A-H,O-Z)
```

Dimension subscripted variables. The dimensions shown permit the cavity to be divided into 100 nodes, with 10 divisions in each direction. Of course, these can be increased by the user if desired.

```
DIMENSION AA(100), CC(100)
DIMENSION VX(100,100), VY(100,100), VZ(100,100)
DIMENSION ZTB(10), DZ(10)
```

Establish COMMON for subroutines called.

```
COMMON /NODE1/ FN1,FN2,N1,N2,M1,M2
COMMON /NODE2/ FNTH,PI,NTH,NN1,NN2,MM2
```

Read in and write out the cavity dimensions and other constants.

```
READ(5,97) A,B,C,N1,N2,NTH,NSHOTN,ABS,REFR,DT,ELIMIT
WRITE(6,96) A,B,C,N1,N2,NTH,NSHOTN,ABS,REFR,DT,ELIMIT
```

Establish the values of the physical constants.

```
COND = 406.64D0
ALPHA = C/(B-A)
SPEC = 234.04D0
THCK = 2.0D-1
PI = 3.1415926D0
DENS = 10524.15D0
```

The cavity is subdivided into $NTH \cdot NN2$ elements, or nodes, in this portion of the program.

```

NARSZ = 100
FNTH = NTH
FN1 = N1
FN2 = N2
NN1 = N1
NN2 = NN1 + N2
M1 = NTH*NN1
M2 = NTH*NN2
MM2 = M2 + 1
I = 0

```

In the "10" DO-loop we compute temporary z-coordinates and the r-coordinate of each node. We use computed GO TO statements to direct the calculation to the proper relations depending on which segment of the cavity is being considered (as determined by the value of "J").

```

DO 10 J = 1, NN2

```

Calculation of the node coordinates, area, thermal capacity and conductances to surrounding nodes is different depending on the surface on which the node is located. Thus we must branch to the appropriate segment of the program.

If the node lies on the barrel, branch to the barrel calculation segment.

```

IF(J.LE.NN1) GOTO 11

```

If the node lies on the cone, branch to the "cone" calculation segment.

```

GOTO 12

```

The node is on the barrel; thus we compute the node coordinates, area, and half-conductances for the barrel.

```
11 CONTINUE
  FJ = J
  TZ = A*(2.0D0*FJ - 1.0D0)/(2.0D0*FN1)
  IF(J.GT.NN2) GOTO 900
  DZ(J) = A/FN1
  ZTB(J) = TZ - DZ(J)/2.0D0
900 CONTINUE
  R = C
  AREA = 2.0D-6*PI*C*A/(FN1*FNTH)
  GK1 = A*COND*THCK*1.0D-3*FNTH/(PI*R*FN1)
  GK2 = COND*PI*R*THCK*4.0D-3/(FNTH*A)
  GK3 = GK1
  GK4 = GK2
  IF(J.EQ.1) GK2 = 0.0D0
  GOTO 17
```

The node is on the cone; thus we compute the coordinates, area, and half-conductances for the cone.

```

12 CONTINUE
  FJ = J
  TZ = 2*(FN1 + FN2 - FJ) + 1.0D0
  TZ = (A-B)*DSQRT(TZ)/DSQRT(2.0D0*FN2) + B
  R = ALPHA*(B-TZ)
  AREA = PI*C*1.0D-6*DSQRT(C*C + (B-A)*(B-A))
  AREA = AREA/(FNTH*FN2)
  TZ1 = 2*(FN1 + FN2 - FJ)
  TZ1 = (A - B)*DSQRT(TZ1)/DSQRT(2.0D0*FN2)
  TZ1 = TZ1 + B
  IF(J.EQ.NN2) GOTO 800
  TZ2 = 2*(FN1 + FN2 - FJ) - 2.0D0
  TZ2 = (A - B)*DSQRT(TZ2)/DSQRT(2.0D0*FN2)
  TZ2 = TZ2 + B
800 TZB2 = (A - B)*DSQRT((FN1+FN2-FJ)/FN2) + B
  TZB1 = (A - B)*DSQRT((FN1+FN2-FJ+1.0D0)/FN2)
  TZB1 = TZB1 + B
  IF(J.GT.NN2) GOTO 901
  ZTB(J) = TZB1
  DZ(J) = TZB2 - TZB1
901 CONTINUE
802 GK1 = (TZB2-TZB1)*(TZB2-TZB1)
  GK1 = GK1 + ALPHA*ALPHA*(TZB2-TZB1)*(TZB2-TZB1)
  GK1 = COND*THCK*FNTH*DSQRT(GK1)*1.0D-3/(PI*R)
  GK2 = (TZ-TZB1)*(TZ-TZB1)
  GK2 = GK2 + ALPHA*ALPHA*(TZ-TZB1)*(TZ-TZB1)
  GK2 = COND*PI*THCK*2.0D-3*ALPHA*(B-TZB1)
  GK2 = GK2/(DSQRT(GK2)*FNTH)
  GK3 = GK1
  IF(J.NE.NN2) GOTO 803
  GK4 = 0.0D0
  GOTO 17
803 GK4 = (TZB2-TZ)*(TZB2-TZ)
  GK4 = GK4 + ALPHA*ALPHA*(TZB2-TZ)*(TZB2-TZ)
  GK4 = COND*PI*THCK*2.0D-3*ALPHA*(TZB2-TZ)
  GK4 = GK4/(DSQRT(GK4)*FNTH)
  GOTO 17

```

This segment of the program is common to both parts of the cavity, barrel and cone. In this section, the coordinates, half-conductances, and areas are stored, and the thermal capacities are computed and stored.

```

17 CONTINUE
  DO 10 K = 1, NTH
    FK = K
    I = I + 1
    VZ(I,I) = TZ
    THETA = 6.283185306D0*FK/FNTH
    VX(I,I) = R*DCOS(THETA)
    VY(I,I) = R*DSIN(THETA)
    IF(I.GT.M2) GOTO 936
    AA(I) = AREA
    CC(I) = AREA*THCK*DENS*SPEC*1.0D-3
936 CONTINUE
    IF(J.GT.NN2) GOTO 10
    ICK = I - 1
    ICK = ICK/NTH
    ICK = I - ICK*NTH
    IF(ICK.EQ.1) GOTO 362
    VX(I,I-1) = GK1
    GOTO 363
362 CONTINUE
    VX(I+NTH-1,I) = GK1
363 CONTINUE
    IF(I.LE.NTH) GOTO 364
    VX(I,I-NTH) = GK2
364 CONTINUE
    IF(ICK.EQ.NTH) GOTO 365
    VY(I+1,I) = GK3
    GOTO 366
365 CONTINUE
    VY(I,I-NTH+1) = GK3
366 CONTINUE
    IF(I.GT.(M2-NTH)) GOTO 10
    VY(I+NTH,I) = GK4
10 CONTINUE

```

(Note: The conductance between node I and adjacent node J is given by $VX(I,J)*VY(I,J)/(VX(I,J) + VY(I,J))$, where $I > J$; that is, series conductances add like parallel resistances.)

Write out results.

```

DO 201 I = 1, M2
  WRITE(10,2) I, AA(I), CC(I)
201 CONTINUE

```

Create a trip card image.

```

ISTAR = 0
AASTAR = 0.0D0
CCSTAR = 0.0D0
WRITE(10,2) ISTAR, AASTAR, CCSTAR

```

Continue writing.

```

DO 202 I = 1, M2
DO 202 J = 1, M2
WRITE(15,1) I, J, VX(I,J), VY(I,J), VZ(I,J)
202 CONTINUE

```

Create another trip card image.

```

I = 0
J = 0
VDUMMY = 0.0D0
WRITE(15,1) I, J, VDUMMY, VDUMMY, VDUMMY

```

And write some more.

```

DO 203 J = 1, NN2
WRITE(15,2) J, ZTB(J), DZ(J)
203 CONTINUE

```

And one last trip card image.

```

J = 0.0D0
ZDUMMY = 0.0D0
WRITE(15,2) J, ZDUMMY, ZDUMMY

```

Subroutine POSMTX is a subroutine which writes out the node position matrix on the line printer so that the values can be verified by the user. Thus, it is optional and does not have to be called.

```

CALL POSMTX(VX,VY,VZ,NARSZ)

```

Format statements:

```

97 FORMAT(F5.2,F5.2,F5.2,I2,I2,I2,I6,F3.1,F3.1,F4.1,
&      E8.2,4I1)
96 FORMAT(5X,'THE SYSTEM VARIABLES ARE AS ',
&      'FOLLOWS:' ,//10X,'A = ',F5.2 , ' MM',
&      /,10X,'B = ',F5.2,' MM',/,10X,'C = ',
&      F5.2,' MM',/,10X,'N1 = ',I2,/,10X,'N2 '
&      ,'= ',I2,/,10X,'NTH =',I2,/,10X,'NSHOTN = ',
&      I6,/,10X,'ABS = ',F3.1,/,10X,'REFR = '
&      ,F3.1,/,10X,'DT = ',F4.1,/,10X,'ELIMIT = '
&      ,E8.2///)
1 FORMAT(5X, 2I5, 3D14.5)
2 FORMAT(5X, I5, 2D14.5)

```

End of Program TRW SHAPE.

```

STOP
END

```


1.3. Subroutine 'POSMTX'

This subroutine rearranges the node coordinate results and writes them out with the appropriate labels so that they can be checked by the user. Use of this subroutine is optional.

```

SUBROUTINE POSMTX(VX,VY,VZ,NARSZ)
  IMPLICIT REAL*8 (A-H,O-Z)
  DIMENSION VX(NARSZ,NARSZ),VY(NARSZ,NARSZ)
  DIMENSION VZ(NARSZ,NARSZ)
  COMMON /NODE1/ FN1,FN2,N1,N2,M1,M2
  COMMON /NODE2/ FNTH,PI,NTH,NN1,NN2,MM2
  WRITE(6,1)
  DO 100 I = 1, M2
    R = DSQRT(VX(I,I)*VX(I,I) + VY(I,I)*VY(I,I))
    PHI = DARCOS(VX(I,I)/R)
    IF(VY(I,I).LT.0.0D0) PHI = 2.0D0*PI - PHI
    WRITE(6,2) I, VX(I,I), VY(I,I), VZ(I,I), R , PHI
100 CONTINUE
  RETURN
  1 FORMAT('1',5X'NODE',T20,'X-COORDINATE',T40,
    &      'Y-COORDINATE',T60,'Z-COORDINATE',T83,
    &      'RADIUS',T103,'ANGLE'//)
  2 FORMAT(6X,I3,T18,D14.5,T38,D14.5,T58,D14.5,
    &      T80,D14.5,T100,D14.5)

```

End of Subroutine 'POSMTX'.

END

2. TRW NODES

This program uses Monte Carlo techniques to compute the distribution factors for the radiation emitted from the cavity nodes for a TRW-type cavity radiometer. The "photon" specular reflections are computed "exactly". Program 'TRW SHAPE' must be executed before this program is submitted.

This program requires one previously generated data file area on system disk. The file should be named A51961.DATA12 and should be of sufficient size to hold $(M2+1)*NN2 + 1$ card images. This file will be written to as unit 12.

2.1. Nomenclature

DF	Matrix of distribution factors. Must be dimensioned $M2 + 1$. DF is written out for each source node.
XPOS	Matrix of node x-coordinates. Must be dimensioned $M2$.
YPOS	Matrix of node y-coordinates. Must be dimensioned $M2$.
ZPOS	Matrix of node z-coordinates. Must be dimensioned $M2$.
ZTB	Matrix of node level upper boundary z-coordinates. Must be dimensioned $NN2$.
DZ	Matrix of node level z-coordinate widths. Must be dimensioned $NN2$.
Z	Matrix of random numbers generated by 'GGUBS', a subroutine called by Subroutine 'RAND'.
A	Distance from plane of precision aperture to junction of barrel and cone (mm).
B	Distance from plane of precision aperture to apex of cone (mm).
C	Radius of barrel (mm).

(Note: For the purpose of stating these dimensions it is assumed that there is no "space ring" gap between the precision aperture and the cavity opening.)

ALPHA	Slope of cone.
NSHOTS	Number of packets into which energy is divided for the Monte Carlo procedure.
RSHOTS	Real value of nshots.
M2	Total number of cavity nodes.
NTH	Number of circumferential divisions.
PI	Ratio of circumference of circle to its diameter.
MM2	Index of aperture.
SEED	Seed number needed for random number generator.
ABS	Absorptivity.
REFR	Ratio of specular component of reflectivity to diffuse + specular (= total) reflectivity.
TXI	X-coordinate of the source position.
TYI	Y-coordinate of the source position.
TZI	Z-coordinate of the source position.
TXIOLD	X-coordinate of the previous source position.
TYIOLD	Y-coordinate of the previous source position.
TZIOLD	Z-coordinate of the previous source position.
TXJ	X-coordinate of the receiving position.
TYJ	Y-coordinate of the receiving position.
TZJ	Z-coordinate of the receiving position.

(Note: All of the above coordinates (TXI, TYI, ..., TZJ) advance with each reflection. Initially TXI, TYI, and TZI are used in a test to determine the first trip through position (TXI, TYI, TZI) by a given "photon".)

NSD Test variable whose value is set to "1" when a specular reflection occurs and to "2" when a diffuse reflection occurs.

2.2. Description of Program

IBM computers require double precision to execute this program. Other compilers may not need this first step:

```
IMPLICIT REAL*8 (A-H, O-Z)
REAL*4 RN
```

Dimension subscripted variables. The subscript range shown below permits the cavity to be divided into up to 100 nodes, with up to 10 divisions in each direction.

```
DIMENSION DF(101), ZTB(10), DZ(10)
DIMENSION XPOS(100), YPOS(100), ZPOS(100)
```

Establish variables common to subroutines called.

```
COMMON /NODE1/ FN1, FN2, N1, N2, NN1
COMMON /NODE2/ FNTH, NTH, NN2, M1, M2, MM2
COMMON /GEOM/ A, B, C, ALPHA, PI
```

Read in and write out cavity dimensions and property values.

```
READ(5, 97) A, B, C, N1, N2, NTH, NSHOTS, ABS, REFR, DT, ELIMIT
WRITE(6, 96) A, B, C, N1, N2, NTH, NSHOTS, ABS, REFR, DT, ELIMIT
```

Establish values of constants.

```
ALPHA = C/(B-A)
PI = 3.141592D0
FNTH = NTH
FN1 = N1
FN2 = N2
NN1 = N1
NN2 = NN1 + N2
M1 = NN1*NTH
M2 = NN2*NTH
MM2 = M2 + 1
RSHOTS = NSHOTS
SEED = 1234567.D0
RN = -1.0D0
```

Read in values of x-, y- and z-positions of nodes; store as XPOS, YPOS and ZPOS, respectively.

```
11 CONTINUE
  READ(15, 1) I, J, XX, YY, ZZ
  IF(I.NE.J) GOTO 11
  IF(I.EQ.0) GOTO 12
  XPOS(J) = XX
  YPOS(J) = YY
  ZPOS(J) = ZZ
  GOTO 11
12 CONTINUE
```

Read in values of ZTB and DZ.

```

899 READ(15,4) J, TZBDUM, DZDUM
    IF(J.EQ.0) GOTO 900
    ZTB(J) = TZBDUM
    DZ(J) = DZDUM
    GOTO 899
900 CONTINUE

```

In the "20" DO-loop we loop through the first nodes of each ring of cavity nodes, treating each as a diffuse source. The distribution factors not computed directly in this loop can be determined when needed from symmetry.

```

I = MM2
DO 20 J = 1, M2, NTH

```

Zero the 'DF' vector.

```

DO 9 I1 = 1, MM2
    DF(I1) = 0.0D0
9 CONTINUE

```

In the "10" DO-loop we allow node J to emit NSHOTS volleys of energy packets and follow them through the cavity until they are either absorbed or leave the cavity through the aperture.

```

DO 10 ISHOT = 1, NSHOTS
    CALL RAND (SEED, IN, RN)
    JL = J - 1
    JL = JL/NTH + 1
    TZI = ZTB(JL) + RN*DZ(JL)
    TZIOLD = TZI
    TZJ = TZI
    CALL RAND(SEED, IN, RN)
    PHII = (2.0D0*RN + 1.0D0)*PI/FNTH
    IF(J.GT.M1) GOTO 902
    R = C
    GOTO 903
902 R = ALPHA*(B - TZI)
903 CONTINUE
    TXI = R*DCOS(PHII)
    TXIOLD = TXI
    TXJ = TXI
    TYI = R*DSIN(PHII)
    TYIOLD = TYI
    TYJ = TYI

```

We return to this junction ("13 CONTINUE") each time a reflection occurs.

```
13 CONTINUE
```

If this is the first time this volley has passed through J, the volley striking node J is "emitted" diffusely.

```
IF((TXI.EQ.TXIOLD).AND.(TYI.EQ.TYIOLD)
&      .AND.(TZI.EQ.TZIOLD)) GOTO 137
```

Test to see if the energy packet is absorbed or reflected. If the random number RN is less than or equal to ABS, the absorptivity, the packet is absorbed; otherwise it is reflected.

```
CALL RAND(SEED, IN, RN)
IF(RN.LE.ABS) GOTO 16
```

Test to see if reflected energy packet is specularly reflected or diffusely reflected. If the random number RN is less than or equal to REFR, the reflectivity ratio, the packet is specularly reflected; otherwise it is diffusely reflected.

```
CALL RAND(SEED, IN, RN)
NSD = 1
IF(RN.LE.REFR) GOTO 17
```

Randomly select the direction of the diffuse reflection.

```
137 CONTINUE
  CALL RAND(SEED, IN, RN)
  IF(RN.EQ.1.0D0) GOTO 14
  THETA = ARSIN(SQRT(RN))
  GOTO 15
14 CONTINUE
  THETA = 0.0D0
15 CONTINUE
  CALL RAND(SEED, IN, RN)
  PHI = 2.0D0*PI*RN
  NSD = 2
  GOTO 17
115 CONTINUE
```

Compute the x-, y-, and z-components of the unit vector in the direction (THETA, PHI) with respect to the central coordinate system.

```
CALL VECTOR(UNX, UNY, UNZ, PHI, THETA, VOX, VOY, VOZ)
```

Call the search subroutine which identifies the receiving point of the diffusely reflected photon.

```
CALL SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)
```

If the z-coordinate of the receiving position is negative, the photon has escaped.

```
IF(TZJ.LT.0.0D0) GOTO 24
```

The photon is reflected to the next position whose coordinates are (TXJ, TYJ, TZJ).

```
GOTO 25
```

The photon is absorbed. Call Subroutine NODE which determines the absorbing node.

```
16 CONTINUE
  CALL NODE(ZTB, TXJ, TYJ, TZJ, J1, JJ1, JJ2)
  DF(J1) = DF(J1) + 1.0D0
  GOTO 10
17 CONTINUE
```

Determine the inward-directed unit normal vector of the source position. This depends on the shape of the surface.

```
IF(TZI.LE.A) GOTO 19
```

The reflecting surface is the cone.

```

18 CONTINUE
  UN = ALPHA*ALPHA*ALPHA*ALPHA
  UN = UN*(B - TZI)*(B - TZI)
  UN = UN + TXI*TXI + TYI*TYI
  UN = DSQRT(UN)
  UNX = - TXI/UN
  UNY = - TYI/UN
  UNZ = - ALPHA*ALPHA*(B - TZI)/UN
  GOTO 23

```

The reflecting surface is the barrel.

```

19 CONTINUE
  UNX = - TXI/C
  UNY = - TYI/C
  UNZ = 0.0D0
23 CONTINUE

```

If reflection is diffuse return to the diffuse sequence;
otherwise continue in the specular sequence.

```

IF(NSD.EQ.2) GOTO 115

```

Compute the "ideal" vector of the reflected packet.

```

VDOT = (TXI-TXIOLD)*UNX
VDOT = VDOT + (TYI-TYIOLD)*UNY
VDOT = VDOT + (TZI-TZIOLD)*UNZ
VOX = TXI - TXIOLD - 2.0D0*VDOT*UNX
VOY = TYI - TYIOLD - 2.0D0*VDOT*UNY
VOZ = TZI - TZIOLD - 2.0D0*VDOT*UNZ
VMAG = DSQRT(VOX*VOX + VOY*VOY + VOZ*VOZ)
VOX = VOX/VMAG
VOY = VOY/VMAG
VOZ = VOZ/VMAG

```

Call the search subroutine which identifies the receiving
point of the "exactly" reflected photon.

```

CALL SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)

```


If the z-coordinate of the receiving position is negative, the photon has escaped.

```
IF(TZJ.LT.0.0D0) GOTO 24
```

The photon is reflected to the next position whose coordinates are (TXJ, TYJ, TZJ).

```
GOTO 25
```

The packet has left the cavity through the aperture. Record and fire another volley.

```
24 CONTINUE
  DF(MM2) = DF(MM2) + 1.0D0
  GOTO 10
```

A reflection to another surface has occurred. Thus, increment TXI, TYI, TZI, TXIOLD, TYIOLD, TZIOLD, TXJ, TYJ, and TZJ.

```
25 CONTINUE
  TXIOLD = TXI
  TYIOLD = TYI
  TZIOLD = TZI
  TXI = TXJ
  TYI = TYJ
  TZI = TZJ
```

Loop back to test sequence to determine disposition of photon when it arrives at this new surface.

```
GOTO 13
10 CONTINUE
```

Write vector of distribution factors corresponding to node J.

```

      JJ = (J + NTH - 1)/NTH
      DO 30 JK = 1, MM2
      DF(JK) = DF(JK)/RSHOTS
      WRITE(12,3) JJ, JK, DF(JK)
30 CONTINUE
20 CONTINUE

```

Generate a "trip" card image at the end of the file.

```

      JJ = 0
      JK = 0
      DFDUM = 0.0D0
      WRITE(12,3) JJ, JK, DFDUM

```

Format statements:

```

1 FORMAT(5X, 2I5, 3D14.5)
2 FORMAT(8X, I3)
3 FORMAT(5X, 2I5, D14.5)
4 FORMAT(5X, I5, 2D14.5)
96 FORMAT(5X, 'THE SYSTEM VARIABLES ARE AS ',
&      'FOLLOWS:' ,//10X, 'A = ', F5.2 , ' MM',
&      /, 10X, 'B = ', F5.2, ' MM' ,/, 10X, 'C = ',
&      F5.2, ' MM' ,/, 10X, 'N1 = ', I2, /, 10X, 'N2 '
&      , '= ', I2, /, 10X, 'NTH = ', I2, /, 10X, 'NSHOTN = '
&      , I6, /, 10X, 'ABS = ', F3.1, /, 10X, 'REFR = '
&      , F3.1, /, 10X, 'DT = ', F4.1, /, 10X, 'ELIMIT = '
&      , E8.2///)
97 FORMAT(F5.2, F5.2, F5.2, I2, I2, I2, I6, F3.1, F3.1, F4.1,
&      E8.2, 4I1)

```

End of Program 'TRW NODES'.

```

STOP
END

```

2.3. Subroutine SEARCH

This subroutine calculates the magnitude of the vector from the source position to the receiving position of the reflected photon. Since the direction of this vector is known (VOX, VOY, VOZ), the magnitude is all that is needed to calculate the receiving position.

```
SUBROUTINE SEARCH(TXI,TYI,TZI,TXJ,TYJ,TZJ,VOX,VOY,VOZ)
```

When this subroutine is compiled on IBM equipment, double precision must be specified. This step may not be necessary on other processors.

```
IMPLICIT REAL*8 (A-H,O-Z)
```

Establish COMMON variables.

```
COMMON /GEOM/ A,B,C,ALPHA,PI
```

If the source position is on the cone and VOZ is positive, search for the receiving position on the cone. If the z-component of the direction unit vector is zero or positive, the escape check may be skipped. Otherwise, begin the search on the aperture, proceeding to the cylinder and cone as necessary.

```
IF((TZI.GE.A).AND.(VOZ.GE.0.0D0)) GOTO 10
```

Check to see if the photon has escaped.

```
IF(VOZ.GE.0.0D0) GOTO 5
VMAG1 = - TZI/VOZ
TXJ = TXI + VMAG1*VOX
TYJ = TYI + VMAG1*VOY
RADIUS = DSQRT(TXJ*TXJ + TYJ*TYJ)
IF(RADIUS.GE.C) GOTO 5
TZJ = - 1.0D0
RETURN
```

Begin the search for the receiving position on the cylinder.

```

5 A1 = VOX*VOX + VOY*VOY
  B1 = 2.0D0*(TXI*VOX + TYI*VOY)
  C1 = TXI*TXI + TYI*TYI - C*C
  VMAG2 = (-B1 + DSQRT(B1*B1 - 4.0D0*A1*C1))
  VMAG2 = VMAG2/(2.0D0*A1)
  VMAG3 = (-B1 - DSQRT(B1*B1 - 4.0D0*A1*C1))
  VMAG3 = VMAG3/(2.0D0*A1)
  VMAGB = VMAG2
  IF(VMAGB.LE.0.0D0) VMAGB = VMAG3
  TXJ = TXI + VMAGB*VOX
  TYJ = TYI + VMAGB*VOY
  TZJ = TZI + VMAGB*VOZ

```

Check to see if the receiving position calculated is on the barrel. If not, continue to search on the cone.

```

IF(TZJ.LE.A) RETURN

```

The receiving position is on the cone.

```

10 A2 = VOX*VOX+VOY*VOY-ALPHA*ALPHA*VOZ*VOZ
  B2 = ALPHA*ALPHA*VOZ*(B-TZI)
  B2 = B2 + TXI*VOX + TYI*VOY
  B2 = B2*2.0D0
  C2 = 2.0D0*B*TZI - TZI*TZI - B*B
  C2 = C2*ALPHA*ALPHA
  C2 = C2 + TXI*TXI + TYI*TYI
  VMAG4 = - B2 + DSQRT(B2*B2 - 4.0D0*A2*C2)
  VMAG4 = VMAG4/(2.0D0*A2)
  VMAG5 = - B2 - DSQRT(B2*B2 - 4.0D0*A2*C2)
  VMAG5 = VMAG5/(2.0D0*A2)
  VMAGC = VMAG4
  IF(VMAGC.LE.0.0D0) VMAGC = VMAG5

```

The equation for the magnitude is quadratic, giving two solutions. The zero or negative root is ignored.

Calculate the receiving position on the cone.

```

TXJ = TXI + VMAGC*VOX
TYJ = TYI + VMAGC*VOY
TZJ = TZI + VMAGC*VOZ
RETURN
END

```

2.4. Subroutine VECTOR

This subroutine converts the randomly selected diffuse angle with respect to the source position into the central coordinate system.

2.4.1. Nomenclature

UNX	X-component of inward-directed unit normal vector of source position.
UNY	Y-component of inward-directed unit normal vector of source position.
UNZ	Z-component of inward-directed unit normal vector of source position.
UTX	X-component of unit tangent vector on source position.
UTY	Y-component of unit tangent vector on source position.
UTZ	Z-component of unit tangent vector on source position.
USX	X-component of the mutually orthogonal vector.
USY	Y-component of the mutually orthogonal vector.
USZ	Z-component of the mutually orthogonal vector.
THETA	Angle of diffuse vector with respect to normal vector.
PHI	Angle of diffuse vector with respect to tangent vector.
VOX	X-component of diffuse vector with respect to central coordinate system.
VOY	Y-component of diffuse vector with respect to central coordinate system.
VOZ	Z-component of diffuse vector with respect to central coordinate system.

2.4.2. Subroutine Description

SUBROUTINE VECTOR(UNX, UNY, UNZ, PHI, THETA, VOX, VOY, VOZ)

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary with other compilers.

IMPLICIT REAL*8 (A-H, O-Z)

Compute the x-, y- and z-components of the unit tangent vector, UT, and the mutually orthogonal vector, US, on the source position.

If the surface is cylindrical, go to 11.

```
IF(UNZ.EQ.0.0D0) GOTO 11
```

Calculations for the general surface.

```
UTX = UNX*DSQRT(UNZ*UNZ/(1.0D0 - UNZ*UNZ))
UTY = UNY*DSQRT(UNZ*UNZ/(1.0D0 - UNZ*UNZ))
UTZ = DSQRT(1.0D0 - UNZ*UNZ)
GOTO 13
```

Calculations for the cylindrical surfaces.

```
11. CONTINUE
   UTX = 0.0D0
   UTY = 0.0D0
   UTZ = 1.0D0
13. CONTINUE
```

Compute the x-, y- and z-components of the unit vector on the source position which is mutually orthogonal to UN and UT.

```
USX = UNY*UTZ - UNZ*UTY
USY = UNZ*UTX - UNX*UTZ
USZ = 0.0D0
```

Compute components of the vector at angle (THETA, PHI) with respect to the UN,UT,US-coordinate system and express the result in the I,J,K-coordinate system.

```
SINTH = DSIN(THETA)
COSTH = DCOS(THETA)
SINPH = DSIN(PHI)
COSPH = DCOS(PHI)
VOX = SINTH*COSPH*UTX + SINTH*SINPH*USX
VOX = VOX + COSTH*UNX
VOY = SINTH*COSPH*UTY + SINTH*SINPH*USY
VOY = VOY + COSTH*UNY
VOZ = SINTH*COSPH*UTZ + SINTH*SINPH*USZ
VOZ = VOZ + COSTH*UNZ
```

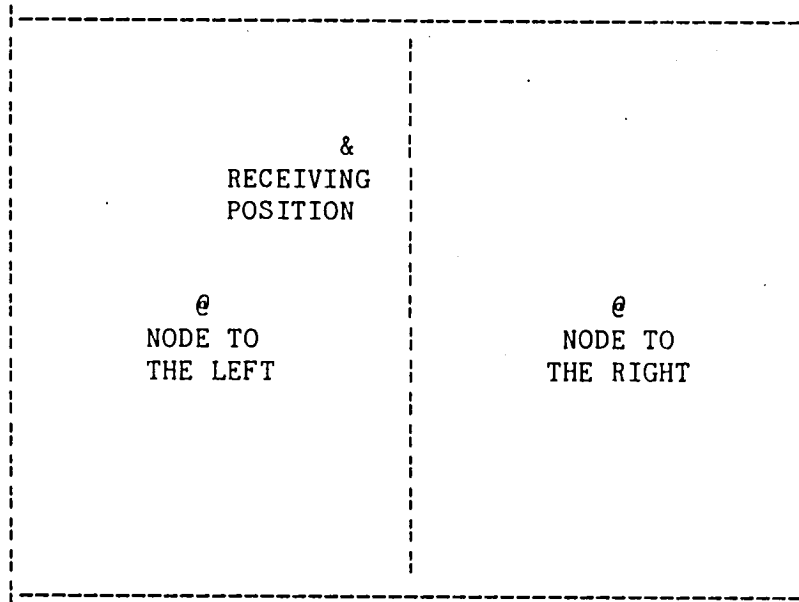
End of Subroutine 'VECTOR'.

RETURN END

2.5. Subroutine NODE

This subroutine identifies the node which absorbs a given photon.

2.5.1. Nomenclature



PHI	Circumferential angle of the receiving position.
PHLE	Circumferential angle of the left edge of the node to the right.
NRA	Number of the node ring "above" the receiving position.
NRB	Number of the node ring "below" the receiving position.
NLP	Position of the node to the left.
NRP	Position of the node to the right.

2.5.2. Subroutine Description

SUBROUTINE NODE(ZTB, TXJ, TYJ, TZJ, J1, JJ1, JJ2)

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary on other compilers.

```
IMPLICIT REAL*8 (A-H, O-Z)
```

Dimension the subscripted variables.

```
DIMENSION ZTB(NN2)
```

Establish COMMON variables.

```
COMMON /NODE1/ FN1, FN2, FN3, N1, N2, N3, NN1
COMMON /NODE2/ FNTH, NTH, NN2, NN3, M1, M2, M3, MM2
COMMON /GEOM/ A, B, C, ALPHA, PI
```

Determine which surface the receiving position lies on. If the receiving position lies on the cone, jump to the cone calculation segment.

```
IF(TZJ.GT.A) GOTO 11
```

The receiving position is on the cylinder.

```
PHI = DAR OS(TXJ/C)
IF(TYJ.LT.0.0DO) PHI = 2.0DO*PI - PHI
NRA = (1.0DO + 2.0DO*FN1*TZJ/A)/2.0DO
NRB = NRA + 1
NRP = FNTH*PHI/(2.0DO*PI)
IF(NRP.EQ.0) NRP = NTH
NLP = NRP + 1
IF(NLP.EQ.(NTH+1)) NLP = 1
PHILE = (2.0DO*NRP + 1.0DO)*PI/FNTH
IF(NRP.EQ.NTH) PHILE = PHILE - 2.0DO*PI
JJ2 = NRP
IF(PHI.GT.PHILE) JJ2 = NLP
JJ1 = NRA
IF(NRA.NE.0) GOTO 998
JJ1 = NRB
GOTO 999
998 IF(TZJ.GT.ZTB(NRB)) JJ1 = NRB
999 J1 = (JJ1 - 1.0DO)*NTH + JJ2
RETURN
```

The receiving position is on the cone.

```

11 R = DSQRT(TXJ*TXJ + TYJ*TYJ)
   PHI = DAR OS(TXJ/R)
   IF(TYJ.LT.0.0D0) PHI = 2.0D0*PI - PHI
   FNRA = (TZJ-B)*(DSQRT(2.0D0*FN2))/(A-B)
   FNRA = FNRA*((TZJ-B)*(DSQRT(2.0D0*FN2))/(A-B))
   FNRA = FN1 + FN2 - (FNRA-1.0D0)/2.0D0
   NRA = FNRA
   NRB = NRA + 1
   NRP = FNTH*PHI/(2.0D0*PI)
   IF(NRP.EQ.0)NRP = NTH
   NLP = NRP + 1
   IF(NLP.EQ.(NTH+1)) NLP = 1
   PHILE = (2.0D0*NRP + 1.0D0)*PI/FNTH
   IF(NRP.EQ.NTH) PHILE = PHILE - 2.0D0*PI
   JJ2 = NRP
   IF(PHI.GT.PHILE) JJ2 = NLP
   JJ1 = NRA
   IF(NRA.EQ.NN2) GOTO 997
   IF(TZJ.GT.ZTB(NRB)) JJ1 = NRB
997 J1 = (JJ1 - 1.0D0)*NTH + JJ2

```

End of Subroutine 'NODE'.

```

RETURN
END

```

2.6. Subroutine RAND

This subroutine uses library Subroutine GGUBS to obtain random numbers in an efficient way. A sequence of 100 random numbers is generated each time the 100th number from the previous generation is used.

2.6.1. Nomenclature

SEED	A seed number required by GGUBS. It must be specified in the main program as a double precision integer value.
IN	An integer between 1 and 100 inclusive which represents which of the 100 random numbers of a given generation is obtained.
RN	The random number passed back to MAIN.

2.6.2. Subroutine Description

```
SUBROUTINE RAND(SEED, IN, RN)
```

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary with other processors.

```
IMPLICIT REAL*8 (A-H,O-Z)
REAL*4 RN, Z(100)
```

If this is the first call to RAND, RN is equal to the negative value set in MAIN. In this case we must call GGUBS to generate the first generation of 100 random numbers.

```
IF(RN.LT.0.0D0) GOTO 11
```

If this is not the first call to RAND, then increment to the next value in this generation of 100 random numbers.

```
IN = IN + 1
```

If we have not yet used up our allotment of random numbers from this generation, we simply take the next one in the sequence back to MAIN.

```
IF(IN.LE.100) GOTO 13
```

However, if we have used all the members of this generation, we must create a new generation of 100 random numbers.

```
11 CONTINUE  
  IN = 1  
  DSEED = SEED  
  NUMBER = 100  
  CALL GGUBS(DSEED, NUMBER, Z)  
  SEED = DSEED  
13 CONTINUE  
  RN = Z(IN)
```

End of Subroutine 'RAND'.

```
RETURN  
END
```

3. TRW FRONT

This program uses Monte Carlo techniques to compute the distribution factors for the radiation emitted from the 'front end' nodes for a TRW-type cavity radiometer. The specular reflections are computed "exactly".

This program requires one previously generated data file area on system disk. The file should be named A51961.DATA17 and should be of sufficient size to hold $(M2+1)*NN2 + 1$ card images. This file will be written to as unit 17, and will contain the distribution factor matrix after the program has been executed.

3.1. Nomenclature

DF	Matrix of distribution factors. Must be dimensioned $M2 + 1$. DF is written out for each source node.
H1	Distance to floor, or face, of instrument (cm).
H2	Distance to face of F-O-V limiter (cm).
R1	Radius which describes the inside face of the F-O-V limiter; also outside radius of the face of the instrument (cm).
R2	Inside radius of the face of the instrument (cm).
R3	Radius of precision aperture (cm).
F1	Radial distance to the inner edge of the first floor node (cm).
F2	Radial distance to the inner edge of the second floor node (cm).
F3	Radial distance to the inner edge of the third floor node (cm).
FOV1	Axial distance from the floor level ($Z = H1$) to the lower edge of the top F-O-V limiter node (cm).
FOV2	Axial distance from the floor level ($Z = H1$) to the lower edge of the central F-O-V limiter node (cm).
EFOV	Emissivity (= absorptivity = $1 - \text{reflectivity}$) of the face of the F-O-V limiter (-).
RFOV	Ratio of specular to specular + diffuse reflectivity for the face of the F-O-V limiter (-).
EF	Emissivity (= absorptivity = $1 - \text{reflectivity}$) of the face of the instrument (-).
RF	Ratio of specular to specular + diffuse reflectivity for the face of the instrument (-).
NSHOTS	Number of "photons" emitted in Monte Carlo method.
RSHOTS	Real value of "NSHOTS".

(Note: The distances H1, H2, R1, R2, R3, and R4 are all measured with respect to the origin of coordinates (0,0,0) located on the axis in the plane of the precision aperture. See Fig. 1 of the February Progress Report.)

PI	Ratio of circumference to diameter of a circle (-).
NTH	Number of circumferential divisions (set at 4).
FNTH	Floating point version of NTH.
N1	Number of axial divisions on F-O-V limiter (-). (set at 3).
FN1	Floating point version of N1.
N2	Number of radial divisions on floor (set at 4).
FN2	Floating point version of N2.
NN2	Number of axial divisions on F-O-V limiter plus number of radial divisions on floor (-).
NN3	Total number of "front end" nodes divided by NTH.
M1	Total number of nodes on F-O-V limiter (-).
M2	Total number of nodes on the face of the instrument including F-O-V limiter and floor (-).
M3	Total number of nodes (= M1 + M2 + 3*NTH), including the edge of the floor, the top surface of the precision aperture, and the imaginary pie shaped segment in the plane of the F-O-V limiter.
SEED	Seed number needed for the random number generator.
RN	Uniformly distributed random number between 0 and 1 generated by Subroutine RAND.
MM3	Total number of possible receiving nodes, including cavity opening.
TXI	X-coordinate of the source position.
TYI	Y-coordinate of the source position.
TZI	Z-coordinate of the source position.
TXIOLD	X-coordinate of the previous source position.
TYIOLD	Y-coordinate of the previous source position.
TZIOLD	Z-coordinate of the previous source position.
TXJ	X-coordinate of the receiving position.
TYJ	Y-coordinate of the receiving position.
TZJ	Z-coordinate of the receiving position.

3.2. Program Description

When this program is compiled on IBM equipment, double precision must be specified. This step may not be necessary on other processors.

```

IMPLICIT REAL*8 (A-H, O-Z)
REAL*4 RN

```

Dimension the subscripted variable.

```

DIMENSION DF(41)

```

Identify the COMMON variables for the subroutines called.

```
COMMON H1,H2,R1,R2,R3,FOV1,FOV2,F1,F2,F3,PI
```

Read in and write out values of front end dimensions and optical properties.

```
READ(5,97) NSHOTS,H1,H2,R1,R2,R3,F1,F2,F3,FOV1,FOV2,
&          EFOV,RFOV,EF,RF
WRITE(6,96) H1,H2,R1,R2,R3,F1,F2,F3,FOV1,FOV2,
&          EFOV,RFOV,EF,RF,NSHOTS
```

Set values of constants.

```
PI = 3.141592D0
NTH = 4
FNTH = NTH
N1 = 3
FN1 = N1
N2 = 4
FN2 = N2
NN2 = N1 + N2
NN3 = NN2 + 3
M1 = N1*NTH
M2 = NN2*NTH
M3 = NN3*NTH
MM3 = M3 + 1
RSHOTS = NSHOTS
SEED = 1234567.D0
RN = -1.0D0
```

In the "20" DO-loop we loop through the first nodes of each ring of NTH nodes, treating each as a diffuse source. The distribution factors not computed directly in this loop can be determined when needed from symmetry.

```
DO 20 J = 1, NN3
JJ = (J-1)*NTH + 1
```

Zero the "DF" vector.

```
DO 9 I1 = 1, MM3
DF(I1) = 0.0D0
9 CONTINUE
```

In the "10" DO-loop we allow node JJ to emit NSHOTS volleys of energy packets and follow them until they are either absorbed or leave the instrument.

```
DO 10 ISHOT = 1, NSHOTS
```

Locate emission point on node JJ.

```
CALL RAND (SEED, IN, RN)
RN1 = RN
CALL RAND(SEED, IN, RN)
RN2 = RN
GOTO(101,102,103,104,105,106,107,108,109,110), J
101 CONTINUE
```

The source point is on the blackbody, which we model as a black sector in the plane of the F-O-V limiter. Note that this limits the analysis to a reasonably uniform calibration source.

```
ANGLE = RN1*PI/2.0D0
RADIUS = DSQRT(R1*R1 - H2*H2)
RADIUS = RADIUS*DSQRT(RN2)
TXI = RADIUS*DCOS(ANGLE)
TYI = RADIUS*DSIN(ANGLE)
TZI = H2
GOTO 111
```

The source point is on the upper ring of the F-O-V limiter face.

```
102 CONTINUE
ANGLE1 = RN1*PI/2.0D0
RAD1 = DSQRT(R1*R1 - H2*H2)
A = DARSIN(RAD1/R1)
RAD2 = DSQRT(R1*R1 - (H1+FOV1)*(H1+FOV1))
B = DARSIN(RAD2/R1)
ANGLE2 = DARCOS(DCOS(A) - RN2*(DCOS(A)-DCOS(B)))
TXI = R1*DSIN(ANGLE2)*DCOS(ANGLE1)
TYI = R1*DSIN(ANGLE2)*DSIN(ANGLE1)
TZI = R1*DCOS(ANGLE2)
GOTO 111
```


The source point is on the middle ring of the F-O-V limiter face.

```

103 CONTINUE
  ANGLE1 = RN1*PI/2.0D0
  RAD1 = DSQRT(R1*R1 - (H1+FOV1)*(H1+FOV1))
  A = DARSIN(RAD1/R1)
  RAD2 = DSQRT(R1*R1 - (H1+FOV2)*(H1+FOV2))
  B = DARSIN(RAD2/R1)
  ANGLE2 = DARCOS(DCOS(A) - RN2*(DCOS(A)-DCOS(B)))
  TXI = R1*DSIN(ANGLE2)*DCOS(ANGLE1)
  TYI = R1*DSIN(ANGLE2)*DSIN(ANGLE1)
  TZI = R1*DCOS(ANGLE2)
  GOTO 111

```

The source point is on the bottom ring of the F-O-V limiter face.

```

104 CONTINUE
  ANGLE1 = RN1*PI/2.0D0
  RAD1 = DSQRT(R1*R1 - (H1+FOV2)*(H1+FOV2))
  A = DARSIN(RAD1/R1)
  RAD2 = DSQRT(R1*R1 - H1*H1)
  B = DARSIN(RAD2/R1)
  ANGLE2 = DARCOS(DCOS(A) - RN2*(DCOS(A)-DCOS(B)))
  TXI = R1*DSIN(ANGLE2)*DCOS(ANGLE1)
  TYI = R1*DSIN(ANGLE2)*DSIN(ANGLE1)
  TZI = R1*DCOS(ANGLE2)
  GOTO 111

```

The source point is on the first (outer) ring of the floor.

```

105 CONTINUE
  ANGLE = RN1*PI/2.0D0
  RAD2 = DSQRT(R1*R1 - H1*H1)
  RADIUS = DSQRT(F1*F1 + RN2*(RAD2*RAD2 - F1*F1))
  TXI = RADIUS*DCOS(ANGLE)
  TYI = RADIUS*DSIN(ANGLE)
  TZI = H1
  GOTO 111

```

The source point is on the second ring of the floor.

```
106 CONTINUE
  ANGLE = RN1*PI/2.0D0
  RADIUS = DSQRT(F2*F2 + RN2*(F1*F1-F2*F2))
  TXI = RADIUS*DCOS(ANGLE)
  TYI = RADIUS*DSIN(ANGLE)
  TZI = H1
  GOTO 111
```

The source point is on the third ring of the floor.

```
107 CONTINUE
  ANGLE = RN1*PI/2.0D0
  RADIUS = DSQRT(F3*F3 + RN2*(F2*F2-F3*F3))
  TXI = RADIUS*DCOS(ANGLE)
  TYI = RADIUS*DSIN(ANGLE)
  TZI = H1
  GOTO 111
```

The source point is on the fourth (inner) ring of the floor.

```
108 CONTINUE
  ANGLE = RN1*PI/2.0D0
  RADIUS = DSQRT(R2*R2 + RN2*(F3*F3-R2*R2))
  TXI = RADIUS*DCOS(ANGLE)
  TYI = RADIUS*DSIN(ANGLE)
  TZI = H1
  GOTO 111
```

The source point is on the inner edge of the floor.

```
109 CONTINUE
  ANGLE = RN1*PI/2.0D0
  ZLEVEL = RN2*H1
  TXI = R2*DCOS(ANGLE)
  TYI = R2*DSIN(ANGLE)
  TZI = ZLEVEL
  GOTO 111
```

The source point is on the precision aperture.

```
110 CONTINUE
  ANGLE = RN1*PI/2.0D0
  RADIUS = R3*DSQRT(RN2)
  TXI = RADIUS*DCOS(ANGLE)
  TYI = RADIUS*DSIN(ANGLE)
  TZI = 0.0D0
```

A reflection to another surface has occurred; thus, increment TXIOLD, TYIOLD, TZIOLD, TXJ, TYJ, and TZJ.

```
111 CONTINUE
  TXIOLD = TXI
  TYIOLD = TYI
  TZIOLD = TZI
  TXJ = TXI
  TYJ = TYI
  TZJ = TZI
```

(Note: All of the above coordinates advance with each reflection. Initially TXI, TYI, and TZI are used in a test to determine the first trip through position (TXI,TYI,TZI) by a given photon.)

We return to this junction ("13 CONTINUE") each time a reflection occurs.

```
13 CONTINUE
```

If this is the first time this volley has passed through J, the volley striking node J is "emitted" diffusely.

```
IF((TXI.EQ.TXIOLD).AND.(TYI.EQ.TYIOLD)
& .AND.(TZI.EQ.TZIOLD)) GOTO 137
```

Test to see if the energy packet is absorbed or reflected. if the random number RN is less than or equal to ABS, the absorptivity, the packet is absorbed; otherwise it is reflected.

```
ABS = EF
REFR = RF
IF(TZJ.GT.H1) ABS = EFOV
IF(TZJ.GT.H1) REFR = RFOV
CALL RAND(SEED, IN, RN)
IF(RN.LE.ABS) GOTO 16
```

Test to see if the reflected energy packet is specularly reflected or diffusely reflected. If the random number RN is less than or equal to REFR, the reflectivity ratio, the packet is specularly reflected; otherwise it is diffusely reflected.

```
CALL RAND(SEED, IN, RN)
NSD = 1
IF(RN.LE.REFR) GOTO 17
```

Randomly select the direction of the diffuse reflection.

```
137 CONTINUE
CALL RAND(SEED, IN, RN)
IF(RN.EQ.1.0D0) GOTO 14
THETA = ARSIN(SQRT(RN))
GOTO 15
14 CONTINUE
THETA = 0.0D0
15 CONTINUE
CALL RAND(SEED, IN, RN)
PHI = 2.0D0*PI*RN
NSD = 2
GOTO 17
115 CONTINUE
```

Call the vector subroutine which computes the x-, y-, and z-components of the unit vector in direction (THETA, PHI) with respect to the central coordinate system.

```
CALL VECTOR(UNX, UNY, UNZ, PHI, THETA, VOX, VOY, VOZ)
```

Call the SEARCH subroutine which identifies the receiving point of the diffusely reflected photon.

```
CALL SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)
```

If the z-coordinate of the receiving position is H2, the photon has escaped.

```
IF(TZJ.EQ.H2) GOTO 24
```

If the z-coordinate of the receiving position is 0 and the r-coordinate is less than R3, the photon has entered the cavity.

```
RAD = DSQRT(TXJ*TXJ + TYJ*TYJ)
IF(TZJ.EQ.0.0D0.AND.RAD.LT.R3) GOTO 244
```

The photon is reflected to the next position whose coordinates are (TXJ, TYJ, TZJ).

```
GOTO 25
```

The photon is absorbed. Call the NODE subroutine which determines the absorbing node.

```
16 CONTINUE
CALL NODE(TXJ, TYJ, TZJ, J1)
DF(J1) = DF(J1) + 1.0D0
GOTO 10
```

Determine the inward-directed unit normal vector of the source position. This depends on the shape of the source surface.

```
17 CONTINUE
IF(TZI.EQ.H2) GOTO 201
IF(TZI.GT.H1) GOTO 202
IF(TZI.EQ.H1) GOTO 203
IF(TZI.GT.0.0D0) GOTO 204
```

The source point is on the precision aperture.

```
UNX = 0.0D0
UNY = 0.0D0
UNZ = 1.0D0
GOTO 23
```

The source is on the imaginary surface in the plane of F-O-V limiter.

```
201 CONTINUE  
    UNX = 0.0D0  
    UNY = 0.0D0  
    UNZ = -1.0D0  
    GOTO 23
```

The source point is on the F-O-V limiter.

```
202 CONTINUE  
    UNX = -TXI/R1  
    UNY = -TYI/R1  
    UNZ = -TZI/R1  
    GOTO 23
```

The source point is on the floor.

```
203 CONTINUE  
    UNX = 0.0D0  
    UNY = 0.0D0  
    UNZ = 1.0D0  
    GOTO 23
```

The source point is on the inner edge of the floor.

```
204 CONTINUE  
    UNX = -TXI/R2  
    UNY = -TYI/R2  
    UNZ = 0.0D0
```

If the reflection is diffuse, return to the diffuse sequence; otherwise continue in the specular sequence.

```
23 CONTINUE  
    IF(NSD.EQ.2) GOTO 115
```

Compute the "ideal" vector of the reflected packet.

```

VDOT = (TXI-TXIOLD)*UNX
VDOT = VDOT + (TYI-TYIOLD)*UNY
VDOT = VDOT + (TZI-TZIOLD)*UNZ
VOX = TXI - TXIOLD - 2.0D0*VDOT*UNX
VOY = TYI - TYIOLD - 2.0D0*VDOT*UNY
VOZ = TZI - TZIOLD - 2.0D0*VDOT*UNZ
VMAG = DSQRT(VOX*VOX + VOY*VOY + VOZ*VOZ)
VOX = VOX/VMAG
VOY = VOY/VMAG
VOZ = VOZ/VMAG

```

Call the SEARCH subroutine which identifies the receiving point of the "exactly" reflected photon.

```

CALL SEARCH(TXI,TYI,TZI,TXJ,TYJ,TZJ,VOX,VOY,VOZ)

```

If the z-coordinate of the receiving position is H2, the photon has escaped.

```

IF(TZJ.EQ.H2) GOTO 24

```

If the z-coordinate of the receiving position is 0 and the r-coordinate is less than R3, the photon has entered the cavity.

```

RAD = DSQRT(TXJ*TXJ + TYJ*TYJ)
IF(TZJ.EQ.0.0D0.AND.RAD.LT.R3) GOTO 244

```

The photon is reflected to the next position whose coordinates are (TXJ, TYJ, TZJ).

```

GOTO 25

```

The photon has escaped through the F-O-V limiter; thus, find the pie segment through which it has escaped and increment the appropriate DF counter.

```
24 CONTINUE
  IF(TXJ.GE.0.0D0.AND.TYJ.GE.0.0D0) I = 1
  IF(TXJ.LT.0.0D0.AND.TYJ.GE.0.0D0) I = 2
  IF(TXJ.LT.0.0D0.AND.TYJ.LT.0.0D0) I = 3
  IF(TXJ.GE.0.0D0.AND.TYJ.LT.0.0D0) I = 4
  DF(I) = DF(I) + 1.0D0
  GOTO 10
```

The photon has entered the cavity; thus, increment the cavity DF counter.

```
244 CONTINUE
  DF(MM3) = DF(MM3) + 1.0D0
  GOTO 10
```

Reflection to another surface has occurred. Thus, increment TXIOLD, TYIOLD, TZIOLD, TXJ, TYJ, and TZJ.

```
25 CONTINUE
  TXIOLD = TXI
  TYIOLD = TYI
  TZIOLD = TZI
  TXI = TXJ
  TYI = TYJ
  TZI = TZJ
```

Loop back to test sequence to determine disposition of the photon when it arrives at this new surface.

```
GOTO 13
```


Write the vector of distribution factors corresponding to node J.

```

10 CONTINUE
   DO 30 JK = 1, MM3
     DF(JK) = DF(JK)/RSHOTS
30 CONTINUE
   DO 63 I = 2, MM3, 4
     I2 = I + 2
     DF(I) = DF(I) + DF(I2)
     DF(I) = DF(I)/2.0D0
     DF(I2) = DF(I)
63 CONTINUE
   DO 64 JK = 1, MM3
     WRITE(17,3) JJ, JK, DF(JK)
64 CONTINUE

```

Generate "trip" card at end of file.

```

20 CONTINUE
   JJ = 0
   JK = 0
   DFDUM = 0.0D0
   WRITE(17,3) JJ, JK, DFDUM

```

Format statements:

```

1 FORMAT(5X, 2I5, 3D14.5)
2 FORMAT(8X, I3)
3 FORMAT(5X, 2I5, D14.5)
4 FORMAT(5X, I5, 2D14.5)
96 FORMAT(5X, 'THE SYSTEM VARIABLES ARE AS ',
&        'FOLLOWS:' ,//10X, 'H1 = ', F5.2, ' CM',
&        '/', 10X, 'H2 = ', F5.2, ' CM' ,/, 10X, 'R1 = ',
&        F5.2, ' CM' ,/, 10X, 'R2 = ', F5.2, '/', 10X, 'R3 '
&        ', ' = ', F5.2, '/', 10X, 'F1 = ', F5.2, '/', 10X, 'F2 = ', F5.2,
&        '/', 10X, 'F3 = ', F5.2, '/', 10X, 'FOV1 = ', F5.2, '/', 10X,
&        'FOV2 = ', F5.2, '/', 10X, 'EFOV = ', F5.2, '/', 10X,
&        'RFOV = ', F5.2, '/', 10X, 'EF = ', F5.2, '/', 10X, 'RF = ',
&        F5.2, '/', 10X, 'NSHOT = ', I6, //)
97 FORMAT(I6, 14F5.2)

```

End of "TRW FRONT".

```

STOP
END

```

3.3. Subroutine SEARCH

This subroutine calculates the magnitude of the vector from the source position to the receiving position of the reflected photon. Since the direction of this vector is known (VOX, VOY, VOZ), the magnitude is all that is needed to calculate the receiving position.

3.3.1. Subroutine Description

```
SUBROUTINE SEARCH(TXI,TYI,TZI,TXJ,TYJ,TZJ,VOX,VOY,VOZ)
```

When this subroutine is compiled on IBM equipment, double precision must be specified. This step may not be necessary on other processors.

```
IMPLICIT REAL*8 (A-H,O-Z)
```

Establish COMMON variables.

```
COMMON H1,H2,R1,R2,R3,FOV1,FOV2,F1,F2,F3,PI
```

If the source point is at the origin of coordinates, RETURN.

```
IF(TZI.EQ.0.0D0.AND.TXI.EQ.0.0D0) GOTO 27
```

If the source point is on the imaginary F-O-V exit plane surface, go to 11.

```
IF(TZI.EQ.H2) GOTO 11
```

If the source point is on the F-O-V limiter, go to 12.

```
IF(TZI.GT.H1) GOTO 12
```

If the source point is on the floor, go to 13.

```
IF(TZI.EQ.H1) GOTO 13
```

If the source point is on the inner edge of the floor, go to 14.

```
IF(TZI.GT.0.0D0) GOTO 14
```

If the source point is on the precision aperture, go to 15.

```
GOTO 15
```

The source point is on the imaginary F-O-V exit plane surface; thus begin search on the F-O-V limiter.

```
11 CONTINUE
VMAG = - (H2-H1)/VOZ
TXJ = TXI + VMAG*VOX
TYJ = TYI + VMAG*VOY
TZJ = H1
RADIUS = DSQRT(TXJ*TXJ + TYJ*TYJ)
```

If 'RADIUS' is less than 'R1', the packet has missed the F-O-V limiter, in which case we go to 16.

```
IF(RADIUS.LT.R1) GOTO 16
```

The packet has intercepted the F-O-V limiter.

```
20 CONTINUE
A = VOX*TXI + VOY*TYI + VOZ*TZI
B = TXI*TXI + TYI*TYI + TZI*TZI - R1*R1
VMAG = - A + DSQRT(A*A-B)
TXJ = TXI + VMAG*VOX
TYJ = TYI + VMAG*VOY
TZJ = TZI + VMAG*VOZ
RETURN
```

The packet has missed the F-O-V limiter. Check to see if it has hit the floor, in which case RETURN.

```
16 CONTINUE
   IF(RADIUS.GT.R2) RETURN
```

The packet has missed the floor; check to see if it has hit the edge of the floor.

```
VMAG = - H2/VOZ
21 CONTINUE
   TXJ = TXI + VMAG*VOX
   TYJ = TYI + VMAG*VOY
   TZJ = 0.0D0
   RADIUS = DSQRT(TXJ*TXJ + TYJ*TYJ)
```

If 'RADIUS' is less than 'R2', the packet has missed the edge of the floor, in which case go to 17.

```
IF(RADIUS.LT.R2) GOTO 17
```

The packet has hit the edge of the floor.

```
25 CONTINUE
   A = VOX*VOX + VOY*VOY
   B = TXI*VOX + TYI*VOY
   C = TXI*TXI + TYI*TYI - R2*R2
   ARG = B*B - A*C
   VMAG = (-B + DSQRT(ARG))/A
   TXJ = TXI + VMAG*VOX
   TYJ = TYI + VMAG*VOY
   TZJ = TZI + VMAG*VOZ
   RETURN
```

The packet has missed the edge of the floor. Check to see if it has hit the precision aperture, in which case RETURN.

```
17 CONTINUE
   IF(RADIUS.GE.R3) RETURN
```

The packet has entered the cavity.

```
TXJ = 0.0D0
TYJ = 0.0D0
TXJ = 0.0D0
RETURN
```

The source point is on the F-O-V limiter; thus begin on the F-O-V limiter exit plane.

```
12 CONTINUE
  IF(VOZ.GT.0.0D0) GOTO 18
```

The source point is not on the F-O-V exit plane; thus begin search on the F-O-V limiter.

```
VMAG = -(TZI-H1)/VOZ
TXJ = TXI + VMAG*VOX
TYJ = TYI + VMAG*VOY
TZJ = H1
RADIUS = DSQRT(TXJ*TXJ + TYJ*TYJ)
```

If 'RADIUS' is less than 'R1', the packet has missed the F-O-V limiter, in which case we continue the search on the floor.

```
IF(RADIUS.LT.R1) GOTO 19
```

The packet has hit the F-O-V limiter. Solve for coordinates.

```
GOTO 20
```

The packet has missed the F-O-V limiter, thus we continue our search on the floor. If 'RADIUS' is greater than 'R2', the packet has hit the floor, in which case we RETURN.

```
19 CONTINUE
  IF(RADIUS.GT.R2) RETURN
```

The packet has missed the floor, in which case we continue our search on the edge of the floor.

```
VMAG = - TZI/VOZ
GOTO 21
```

Search on the F-O-V limiter exit plane.

```
18 CONTINUE
   VMAG = (H2-TZI)/VOZ
22 CONTINUE
   TXJ = TXI + VMAG*VOX
   TYJ = TYI + VMAG*VOY
   TZJ = H2
   RADIUS = DSQRT(TXJ*TXJ + TYJ*TYJ)
   REXIT = DSQRT(R1*R1 - H2*H2)
```

If 'RADIUS' is less than 'REXIT', the packet has escaped, in which case we RETURN.

```
IF(RADIUS.LT.REXIT) RETURN
```

The packet has hit the F-O-V limiter.

```
GOTO 20
```

The source point is on the floor; thus begin search on the imaginary F-O-V exit plane surface.

```
13 CONTINUE
   VMAG = (H2-H1)/VOZ
   GOTO 22
```

The source point is on the inner edge of the floor; thus check to see if packet is emitted upward or downward.

```
14 CONTINUE
   IF(VOZ.LT.0.0D0) GOTO 23
```

The packet has been emitted upward; thus resume the search on the inner edge of the floor.

```

      VMAG = (H1-TZI)/VOZ
26  CONTINUE
      TXJ = TXI + VMAG*VOX
      TYJ = TYI + VMAG*VOY
      RADIUS = DSQRT(TXJ*TXJ + TYJ*TYJ)

```

If 'RADIUS' is less than 'R2', the packet has missed the inner edge of the floor; thus resume search on the F-O-V limiter.

```

      IF(RADIUS.LT.R2) GOTO 24

```

The packet has intercepted the inner edge of the floor; thus go to 25.

```

      GOTO 25

```

The packet has missed the inner edge of the floor; thus resume search on the F-O-V limiter.

```

24  CONTINUE
      VMAG = (H2-TZI)/VOZ
      GOTO 22

```

The packet is emitted downward; thus begin search on the inner edge of the floor.

```

23  CONTINUE
      VMAG = -TZI/VOZ
      TXJ = TXI + VMAG*VOX
      TYJ = TYI + VMAG*VOY
      TZJ = 0.0D0
      RADIUS = DSQRT(TXJ*TXJ + TYJ*TYJ)

```

If 'RADIUS' is greater than 'R2', the packet has hit the inner edge of the floor, in which case go to 25; otherwise, go to 17.

```

      IF(RADIUS.GT.R2) GOTO 25
      GOTO 17

```

The source point is on the precision aperture; thus begin search on the inner edge of the floor.

```

15 CONTINUE
   VMAG = H1/VOZ
   GOTO 26
27 CONTINUE

```

End of Subroutine 'SEARCH'.

```

RETURN
END

```

3.4. Subroutine VECTOR

This subroutine expresses the randomly selected diffuse angle with respect to the source position in terms of the central coordinate system.

3.4.1. Nomenclature

UNX	X-component of inward-directed unit normal vector of source position.
UNY	Y-component of inward-directed unit normal vector of source position.
UNZ	Z-component of inward-directed unit normal vector of source position.
UTX	X-component of unit tangent vector of source position.
UTY	Y-component of unit tangent vector of source position.
UTZ	Z-component of unit tangent vector of source position.
USX	X-component of the mutually orthogonal vector.
USY	Y-component of the mutually orthogonal vector.
USZ	Z-component of the mutually orthogonal vector.
THETA	Angle of diffuse vector with respect to normal vector.
PHI	Angle of diffuse vector with respect to tangent vector.
VOX	X-component of diffuse vector with respect to central coordinate system.
VOY	Y-component of diffuse vector with respect to central coordinate system.
VOZ	Z-component of diffuse vector with respect to central coordinate system.

3.4.2. Subroutine Description

```
SUBROUTINE VECTOR(UNX, UNY, UNZ, PHI, THETA, VOX, VOY, VOZ)
```

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary on other processors.

```
IMPLICIT REAL*8 (A-H, O-Z)
```

Compute the x-, y-, and z-components of the unit tangent vector, UT, and the mutually orthogonal vector, US, on the source position.

If the surface is cylindrical, go to 11.

```
IF(UNZ.EQ.0.0D0) GOTO 11
```

If the surface is plane, go to 12.

```
IF(UNZ.EQ.1.0D0.OR.UNZ.EQ.-1.0D0) GOTO 12
```

Calculations for the spherical surface:

```
ARG = UNZ*UNZ/(1.0D0-UNZ*UNZ)
UTX = UNX*DSQRT(ARG)
UTY = UNY*DSQRT(ARG)
UTZ = DSQRT(1.0D0 - UNZ*UNZ)
GOTO 13
```

Calculations for the cylindrical surface:

```
11 CONTINUE
   UTX = 0.0D0
   UTY = 0.0D0
   UTZ = 1.0D0
   GOTO 13
```

Calculations for plane surfaces:

```
12 CONTINUE
   UTX = 1.0D0
   UTY = 0.0D0
   UTZ = 0.0D0
13 CONTINUE
```

Compute the x-, y-, and z-components of the unit vector on the source position which is mutually orthogonal to UN and UT.

```
USX = UNY*UTZ - UNZ*UTY
USY = UNZ*UTX - UNX*UTZ
USZ = 0.0D0
```

Compute components of the vector at angle (THETA, PHI) with respect to the UN,UT,US-coordinate system and express the result in the I,J,K-coordinate system.

```
SINTH = DSIN(THETA)
COSTH = DCOS(THETA)
SINPH = DSIN(PHI)
COSPH = DCOS(PHI)
VOX = SINTH*COSPH*UTX + SINTH*SINPH*USX
VOX = VOX + COSTH*UNX
VOY = SINTH*COSPH*UTY + SINTH*SINPH*USY
VOY = VOY + COSTH*UNY
VOZ = SINTH*COSPH*UTZ + SINTH*SINPH*USZ
VOZ = VOZ + COSTH*UNZ
```

End of Subroutine 'VECTOR'.

```
RETURN
END
```

3.5. Subroutine NODE

This subroutine identifies the node which absorbs a given photon.

3.5.1. Subroutine Description

```
SUBROUTINE NODE(TXJ, TYJ, TZJ, J1)
```

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary on other processors.

```
IMPLICIT REAL*8 (A-H, O-Z)
```

Identify the COMMON variables.

```
COMMON H1, H2, R1, R2, R3, FOV1, FOV2, F1, F2, F3, PI
```

Check to see if absorption occurred on the F-O-V limiter, in which case go to 11.

```
IF(TZJ.GT.H1) GO TO 11  
RADIUS = DSQRT(TXJ*TXJ + TYJ*TYJ)
```

Check to see if absorption occurred on the floor, in which case go to 12.

```
IF(RADIUS.LT.R1.AND.RADIUS.GE.R2) GO TO 12
```

Check to see if absorption occurred on the edge of the floor, in which case go to 13.

```
IF(TZJ.LT.H1.AND.TZJ.GT.0.0D0) GOTO 13
```

Absorption occurred on the precision aperture; thus determine which node.

```
IF(TXJ.GE.0.0DO.AND.TYJ.GE.0.0DO) J1 = 37
IF(TXJ.LT.0.0DO.AND.TYJ.GE.0.0DO) J1 = 38
IF(TXJ.LT.0.0DO.AND.TYJ.LT.0.0DO) J1 = 39
IF(TXJ.GE.0.0DO.AND.TYJ.LT.0.0DO) J1 = 40
RETURN
```

Absorption occurred on the F-O-V limiter; thus check to see in which ring of nodes the absorption occurred.

```
11 CONTINUE
IF(TZJ.GT.FOV1) GOTO 14
IF(TZJ.GT.FOV2) GOTO 15
```

Absorption occurred in the bottom ring of nodes of the F-O-V limiter; thus identify the node.

```
IF(TXJ.GE.0.0DO.AND.TYJ.GE.0.0DO) J1 = 13
IF(TXJ.LT.0.0DO.AND.TYJ.GE.0.0DO) J1 = 14
IF(TXJ.LT.0.0DO.AND.TYJ.LT.0.0DO) J1 = 15
IF(TXJ.GE.0.0DO.AND.TYJ.LT.0.0DO) J1 = 16
RETURN
```

Absorption occurred in the top ring of nodes of the F-O-V limiter; thus identify the node.

```
14 CONTINUE
IF(TXJ.GE.0.0DO.AND.TYJ.GE.0.0DO) J1 = 5
IF(TXJ.LT.0.0DO.AND.TYJ.GE.0.0DO) J1 = 6
IF(TXJ.LT.0.0DO.AND.TYJ.LT.0.0DO) J1 = 7
IF(TXJ.GE.0.0DO.AND.TYJ.LT.0.0DO) J1 = 8
RETURN
```

Absorption occurred in the middle ring of nodes of the F-O-V limiter; thus identify the node.

```
15 CONTINUE
IF(TXJ.GE.0.0DO.AND.TYJ.GE.0.0DO) J1 = 9
IF(TXJ.LT.0.0DO.AND.TYJ.GE.0.0DO) J1 = 10
IF(TXJ.LT.0.0DO.AND.TYJ.LT.0.0DO) J1 = 11
IF(TXJ.GE.0.0DO.AND.TYJ.LT.0.0DO) J1 = 12
RETURN
```

Absorption occurred on the floor; thus check to see in which ring of nodes the absorption occurred.

```
12 CONTINUE
  IF(RADIUS.GE.F1) GOTO 16
  IF(RADIUS.GT.F2) GOTO 17
  IF(RADIUS.GT.F3) GOTO 18
```

Absorption occurred in the innermost ring of nodes on the floor; thus identify the node.

```
  IF(TXJ.GE.0.0D0.AND.TYJ.GE.0.0D0) J1 = 29
  IF(TXJ.LT.0.0D0.AND.TYJ.GE.0.0D0) J1 = 30
  IF(TXJ.LT.0.0D0.AND.TYJ.LT.0.0D0) J1 = 31
  IF(TXJ.GE.0.0D0.AND.TYJ.LT.0.0D0) J1 = 32
  RETURN
```

Absorption occurred on the outermost ring of nodes on the floor; thus identify the node.

```
16 CONTINUE
  IF(TXJ.GE.0.0D0.AND.TYJ.GE.0.0D0) J1 = 17
  IF(TXJ.LT.0.0D0.AND.TYJ.GE.0.0D0) J1 = 18
  IF(TXJ.LT.0.0D0.AND.TYJ.LT.0.0D0) J1 = 19
  IF(TXJ.GE.0.0D0.AND.TYJ.LT.0.0D0) J1 = 20
  RETURN
```

Absorption occurred on the second ring of nodes from the outer edge of the floor; thus identify the node.

```
17 CONTINUE
  IF(TXJ.GE.0.0D0.AND.TYJ.GE.0.0D0) J1 = 21
  IF(TXJ.LT.0.0D0.AND.TYJ.GE.0.0D0) J1 = 22
  IF(TXJ.LT.0.0D0.AND.TYJ.LT.0.0D0) J1 = 23
  IF(TXJ.GE.0.0D0.AND.TYJ.LT.0.0D0) J1 = 24
  RETURN
```

Absorption occurred on the third ring of nodes from the outer edge of the floor; thus identify the node.

```
18 CONTINUE
  IF(TXJ.GE.0.0D0.AND.TYJ.GE.0.0D0) J1 = 25
  IF(TXJ.LT.0.0D0.AND.TYJ.GE.0.0D0) J1 = 26
  IF(TXJ.LT.0.0D0.AND.TYJ.LT.0.0D0) J1 = 27
  IF(TXJ.GE.0.0D0.AND.TYJ.LT.0.0D0) J1 = 28
  RETURN
```

Absorption occurred on the edge of the floor; thus
identify the node.

```
13 CONTINUE  
  IF(TXJ.GE.0.0D0.AND.TYJ.GE.0.0D0) J1 = 33  
  IF(TXJ.LT.0.0D0.AND.TYJ.GE.0.0D0) J1 = 34  
  IF(TXJ.LT.0.0D0.AND.TYJ.LT.0.0D0) J1 = 35  
  IF(TXJ.GE.0.0D0.AND.TYJ.LT.0.0D0) J1 = 36
```

End of Subroutine 'NODE'.

```
RETURN  
END
```

3.6. Subroutine RAND

This subroutine uses library Subroutine GGUBS to obtain random numbers in an efficient way. A sequence of 100 random numbers is generated each time the 100th number from the previous generation is used.

3.6.1. Nomenclature

SEED	A seed number required by GGUBS. It must be specified in the main program as a double precision integer value.
IN	An integer between 1 and 100 inclusive which represents which of the 100 random numbers of a given generation is obtained.
RN	The random number passed back to MAIN.

3.6.2. Subroutine Description

```
SUBROUTINE RAND(SEED, IN, RN)
```

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary with other processors.

```
IMPLICIT REAL*8 (A-H,O-Z)
REAL*4 RN, Z(100)
```

If this is the first call to RAND, RN is equal to the negative value set in MAIN. In this case we must call GGUBS to generate the first generation of 100 random numbers.

```
IF(RN.LT.0.0D0) GOTO 11
```

If this is not the first call to RAND, then increment to the next value in this generation of 100 random numbers.

```
IN = IN + 1
```

If we have not yet used up our allotment of random numbers from this generation, we simply take the next one in the sequence back to MAIN.

```
IF(IN.LE.100) GOTO 13
```

However, if we have used all the members of this generation, we must create a new generation of 100 random numbers.

```
11 CONTINUE  
  IN = 1  
  DSEED = SEED  
  NUMBER = 100  
  CALL GGUBS(DSEED, NUMBER, Z)  
  SEED = DSEED  
13 CONTINUE  
  RN = Z(IN)
```

End of Subroutine 'RAND'.

```
RETURN  
END
```


<u>Variable</u>	<u>Type</u>	<u>Field</u>	<u>Description</u>
NSHOT	I6	1-6	Number of volleys fired
H1	F5.2	7-11	Distance to floor of instrument
H2	F5.2	12-16	Distance to face of F-O-V limiter
R1	F5.2	17-21	Outside radius of face
R2	F5.2	22-26	Inside radius of face
R3	F5.2	27-31	Radius of precision aperture
F1	F5.2	32-36	Radius to first floor node
F2	F5.2	37-41	Radius to second floor node
F3	F5.2	42-46	Radius to third floor node
FOV1	F5.2	47-51	Distance to top F-O-V limiter node
FOV2	F5.2	52-56	Distance to center limiter node
EFOV	F5.2	57-61	Emissivity of F-O-V limiter face
RFOV	F5.2	62-66	Reflectivity ratio of F-O-V face
EF	F5.2	67-71	Emissivity of floor
RF	F5.2	72-76	Reflectivity ratio of face

Note: Please see the 'NOMENCLATURE' of this section for a more complete description of these variables. Also, see Fig. 1 of the February Progress Report.

4. TRW TEMP

This program computes the transient temperature distribution in the active cavity and body of a TRW-type wide-field-of-view total spectrum radiometer subject to a uniform blackbody irradiance input and arbitrary heat sink and mounting beam temperature variations. It requires as inputs conductance matrices and radiation distribution factor matrices generated by the previous programs.

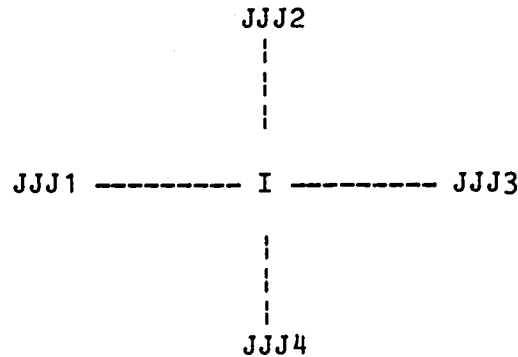
Programs TRW SHAPE, TRW NODES and TRW FRONT must be executed before this program is submitted. Also, an additional on-line disk file must be created and named A51961.DAT14 which contains the body internode conductances and the body node surface areas, emissivities, thermal capacities, volumes and heat sources. The current version of this file, which is user generated, is given in section 6. of this USER'S MANUAL.

4.1. Nomenclature

TO	Matrix of "old" cavity node temperatures (K); i.e., temperatures at beginning of calculation time interval DT. Must be dimensioned M2.
TN	Matrix of "new" cavity node temperatures (K); i.e., temperatures at the end of calculation time interval DT. Must be dimensioned M2.
AA	Matrix of cavity node surface areas (m**2). Must be dimensioned M2.
CC	Matrix of cavity node thermal capacities (W-s/K). Must be dimensioned M2.
DFC(J,I)	Matrix of cavity distribution factors for radiation emitted diffusely from the first node position on node ring J which is absorbed by node I (-). Must be dimensioned NN2 by (M2 + 1).
G(I,J)	Matrix of conductances for conduction path between cavity nodes I and J (W/K). Must be dimensioned M2 by M2.
NBODY	Number of radiometer "body" nodes (= 104).
NBODY2	NBODY + 2 (= 106).
RNBODY	Real version of NBODY.
HST	Heat sink (node "20") temperature (K).
TCLAMP	Mounting beam temperature (K).
HEAT	Proportional to cavity heater power, P (P = 0.32*HEAT mW).
TSOURC	Source equivalent blackbody temperature (K).
GLINK	The conductance of the link between the cavity and the thermal impedance (W/K).
SIG	Stefan-Boltzmann constant (5.6693E-08 W/m**2-K**4).
MBODY	Body node number.

AABODY Surface area of body node 'MBODY' for purposes of calculating radiation heat transfer (cm^2).
 EEBODY Emissivity of body node 'MBODY' (-).
 CCBODY Thermal capacity of body node 'MBODY' ($\text{W-s/cm}^3\text{-K}$).
 VVBODY Volume of body node 'MBODY' (cm^3).
 QQBODY Heat source in body cavity 'MBODY'; may be due to radiation, conduction or electrical heating (W).
 LIMBODY Delimiting index in DO-loop ($= \text{MBODY} + 3$).
 ABODY(I) Surface area of body node I for purposes of calculating radiation exchange (m^2).
 EBODY(I) Emissivity of body node I (-).
 CBODY(I) Thermal capacity of body node I ($\text{W-s/m}^3\text{-K}$).
 VBODY(I) Volume of body node I (m^3).
 QBODY(I) Heat source in body cavity I; may be due to radiation, conduction or electrical heating (W).
 T1BODY(I) Initial temperature of body node I at beginning of solution; must be specified by user by specifying 'HST'. Later set equal to new temperature of node I at end of calculation interval (K).
 TGBODY(I) Initial guess of body node I temperature at the end of a calculation time interval; initially set equal to 'HST', but in subsequent time steps it is updated to previous value (K).
 COND Value of thermal conductance between body nodes on the same level ('MBODY') in the azimuthal direction (W/K).
 GBODY(I,J) Value of thermal conductance between body nodes I and J (W/K).
 DBODY(I,J) Value of radiation distribution factor from body node I to body node J (-).
 MIBODY Index number of body node "level", or ring (-).
 MJBODY Index number of body node "level", or ring (-).
 SOURCE(I) Blackbody source-to-body node I distribution factor (-).
 BODYDF Value of distribution factor, either body node-to-body node or source-to-body node, read from unit 17 (-).
 FTIME Total calculation time; i.e., duration of the problem (s).
 DT Length of cavity calculation time increment; i.e., integration step size (s).
 NDT Number of time increments used in cavity calculation (-).
 DTBODY Length of body calculation time increment (s).
 PI Ratio of circumference to diameter of a circle (-).
 NARSZ Index used in Subroutine GAUSS; represents upper limit of size of matrix that can be inverted by the subroutine; i.e., used as argument of dimension statement in the subroutine. (-).
 M2 Number of cavity nodes (-).
 ELIMIT RMS error limit (%).
 ITRMAX Maximum number of iterations permitted in the cavity temperature distribution calculation loop.

ITRMIN Minimum number of iterations permitted in the
 cavity temperature distribution calculation loop.
 AO Aperture area (mm**2).
 NTH Number of circumferential divisions of cavity (-).
 EC Emissivity (=absorptivity) of the cavity (-).
 EB Emissivity of the back surface of the cavity (-).
 ETIME Elapsed time counter (s).
 QTOT(I) Average rate of energy transfer from the cavity
 to Node '13' of the radiometer body during ten
 successive cavity calculation time increments (W).
 I1 Index indicating level of Node I.
 I2 Index indicating position of Node I in Level I1.
 JJJ1 Index of node "to left" of Node I.
 JJJ2 Index of node "above" Node I.
 JJJ3 Index of Node "to right" of Node I.
 JJJ4 Index of node "below" Node I.



Q1 Energy transfer rate into Node I from cavity nodes
 other than the receiving node itself during
 calculation time interval DT (W).
 J1 Index indicating level of Node J.
 J2 Index indicating position of node J in Level J1.

4.2. Program Description

Double precision must be specified when this program is compiled on IBM equipment. When other processors are used, the following statement may not be necessary.

IMPLICIT REAL*8 (A-H,O-Z)

Dimension subscripted variables.

```
DIMENSION ABODY(104),EBODY(104),CBODY(104),VBODY(104)
DIMENSION QBODY(104),T1BODY(104),TGBODY(104),TUBODY(104)
DIMENSION QTOT(4),SOURCE(45),IANS(100)
DIMENSION DBODY(104,106),GBODY(104,104)
DIMENSION TO(100),TN(100),AA(100),CC(100),RHS(100)
DIMENSION Z(100,100),G(100,100)
DIMENSION DFC(10,101),DFA(4,100,101)
```

Establish COMMON variables.

```
COMMON /EMSGC/ IER
```

Set constants. (Note: This is where each of these numbers must be changed when different cases are to be run.)

```
NBODY = 104
RNBODY = NBODY
HST = 313.0D0
TCLAMP = 326.0D0
HEAT = 400.0D0
TSOURC = 299.36D0
GLINK = 0.04D0
SIG = 5.6693D-08
NARSZ = 100
PI = 3.141592D0
```

Read in and write out system input variables.

```
READ(5,97) A,B,C,N1,N2,NTH,NSHOTN,ABS,REFR,DT,ELIMIT,MAT
WRITE(6,96) A,B,C,N1,N2,NTH,NSHOTN,ABS,REFR,DT,ELIMIT,MAT
```

The following 'WRITE' statement causes a dimensioned sketch of the cavity to be "drawn" by the line printer.

```
WRITE(6,94)
```

Initialize radiometer body internode conductances and distribution factors to zero. This is necessary because the solution strategy considers possible interaction between all possible combinations of any two radiometer body nodes. Later we will read in the nonzero values of these two matrices.

```

      NBODY2 = NBODY + 2
      DO 111 I = 1,NBODY
      DO 111 J = I,NBODY2
      IF(J.LE.NBODY) GBODY(I,J) = 0.0D0
      IF(J.LE.NBODY) GBODY(J,I) = GBODY(I,J)
      DBODY(I,J) = 0.0D0
      IF(J.LE.NBODY) DBODY(J,I) = DBODY(I,J)
111  CONTINUE

```

Read in values of constants and properties; initialize the node temperatures for radiometer body.

```

123  CONTINUE
      READ(14,81) MBODY,AABODY,EEBODY,CCBODY,VVBODY,QQBODY
      IF(MBODY.EQ.0) GOTO 124
      LIMBOD = MBODY + 3
      DO 223 I = MBODY,LIMBOD
      ABODY(I) = AABODY*1.0D-4
      EBODY(I) = EEBODY
      CBODY(I) = CCBODY*1.0D6
      VBODY(I) = VVBODY*1.0D-6
      QBODY(I) = QQBODY
      T1BODY(I) = HST
      TGBODY(I) = HST
223  CONTINUE
      GOTO 123
124  CONTINUE

```

Read in nonzero elements of 'GBODY' matrix.

```
211 CONTINUE
    READ(14,82) MBODY,COND
    IF(MBODY.EQ.0) GOTO 212
    LIMBOD = MBODY + 3
    DO 215 I = MBODY,LIMBOD
        J = I + 1
        IF(I.EQ.LIMBOD) J = MBODY
        GBODY(I,J) = COND
        GBODY(J,I) = COND
215 CONTINUE
    GOTO 211
212 CONTINUE
    READ(14,83) MIBODY,MJBODY,GBODY(MIBODY,MJBODY)
    IF(MIBODY.EQ.0) GOTO 213
    GBODY(MJBODY,MIBODY) = GBODY(MIBODY,MJBODY)
    GOTO 212
213 CONTINUE
```

Initialize source-to-instrument distribution factors to zero.

```
DO 601 I = 1,41
    SOURCE(I) = 0.0D0
601 CONTINUE
```


Read in nonzero values of 'DBODY' matrix.

```

      INN = 0
      INK = 1
403  CONTINUE
      READ(17,92) I, J, BODYDF
      IF(I.EQ.0) GOTO 402
      INN = INN + 1
      IF(I.EQ.1) GOTO 501
      IF(I.LE.13) II = I - 4
      IF(I.EQ.17) II = 17
      IF(I.GE.21.AND.I.LT.33) II = I + 8
      IF(I.GE.33) II = I + 4
      IF(J.LT.5) JJ = 105
      IF(J.GE.5.AND.J.LT.17) JJ = J - 4
      IF(J.GE.17.AND.J.LT.21) JJ = J
      IF(J.GE.21.AND.J.LT.33) JJ = J + 8
      IF(J.GE.33.AND.J.LT.41) JJ = J + 4
      IF(J.EQ.41) JJ = 106
      DBODY(II,JJ) = DBODY(II,JJ) + BODYDF
      GOTO 403
501  CONTINUE
      IF(INN.EQ.5) INK = INK + 4
      IF(INN.EQ.5) INN = 1
      SOURCE(INK) = SOURCE(INK) + BODYDF
      GOTO 403
402  CONTINUE

```

Use symmetry to establish the values of the elements of the 'DBODY' matrix not computed directly.

```

      DO 401 I = 1,104,4
      DO 401 K = 1,104
      I1 = I + 1
      I2 = I + 2
      I3 = I + 3
      L1 = K + 3
      IF(L1.GT.4) L1 = K - 1
      L2 = K + 2
      IF(L2.GT.4) L2 = K - 2
      L3 = K + 1
      IF(L3.GT.4) L3 = K - 3
      DBODY(I1,K) = DBODY(I,L1)
      DBODY(I2,K) = DBODY(I,L2)
      DBODY(I3,K) = DBODY(I,L3)
401  CONTINUE

```

Set upper limit on calculation time; i.e., the duration of the heating or cooling problem being studied. The value used below, 400 s, permits the instrument to come into equilibrium. Also, set number of calculation time intervals.

```
FTIME = 400.0D0
NDT = FTIME/DT
```

Set calculation step size for body temperature calculation. Note that this is ten times the calculation time increment for the cavity calculation.

```
DTBODY = 10.0D0*DT
```

Establish constants and indices.

```
AO = PI*C*C/100.0D0
FNTH = NTH
NN1 = N1
NN2 = NN1 + N2
M1 = NN1*NTH
M2 = NN2*NTH
FM2 = M2
MM2 = M2 + 1
EC = ABS
ITRMAX = 100
ITRMIN = 2
EB = 1.0D0
ETIME = 0.0D0
IER = 0
```

Write out node grid and geometry information.

```
WRITE(6,98) M2,NN2,NTH,NA,NTH,A,B,C
IF(MAT.EQ.0) WRITE(6,93) ELIMIT
```

Read in values of dimensioned constants. Begin by zeroing the G(I,J), Z(I,J) and RHS(I) matrices.

```
DO 19 I = 1, M2
  RHS(I) = 0.0D0
DO 19 J = 1, M2
  G(I,J) = 0.0D0
  Z(I,J) = 0.0D0
19 CONTINUE
```

Note: The node half-conductances must be retrieved from the lower left-hand triangle of the 'V' matrix created by 'TRW SHAPE'.

```

10 CONTINUE
  READ(15,7) I, J, XX, YY, ZZ
  IF((I.GT.M2).OR.(I.EQ.0)) GOTO 11
  IF(I.LE.J) GOTO 10

```

If I or J are either one greater than M2, the values of XX and YY will be zero; that is, there are no conductances stored at those locations. We skip the following two statements in that case.

```

  IF(XX.LE.1.0D-10) GOTO 10
  IF(YY.LE.1.0D-10) GOTO 10
  G(I,J) = 1.0D0/XX + 1.0D0/YY
  G(I,J) = 1.0D0/G(I,J)
  G(J,I) = G(I,J)
  GOTO 10
11 CONTINUE
  READ(10,1) I, AI, CI
  IF(I.EQ.0) GOTO 12
  AA(I) = AI
  CC(I) = CI
  GOTO 11
12 CONTINUE
  READ(11,2) L, M, I, DFADUM
  IF(L.EQ.0) GOTO 13
  DFA(L,M,I) = DFADUM
  GOTO 12
13 CONTINUE
  READ(12,3) I, J, DFCIJ
  IF(I.EQ.0) GOTO 14
  DFC(I,J) = DFCIJ
  GOTO 13
14 CONTINUE

```

Initialize 'old' node temperatures. This establishes the initial conditions for the solution.

```

DO 15 I = 1, M2
  TO(I) = HST
  TN(I) = TO(I)
15 CONTINUE

```

Write out initial cavity temperature distribution.

```

WRITE(6,8) ETIME
DO 301 I5 = NTH, M2 , NTH
  I9 = I5 - NTH + 1
  WRITE(6,6) (TN(I3), I3 = I9, I5)
  WRITE(16,6) (TN(I3), I3 = I9, I5)
301 CONTINUE

```

Zero the 'QTOT' accumulator.

```

DO 987 I = 1,4
  QTOT(I) = 0.0D0
987 CONTINUE

```

We "march in time" by letting (if desired) the heat sink and calibration source temperatures change with each time increment DT. We assume that DT is sufficiently small that each incremental change in these temperatures is small.

The source and heat sink temperatures are changed in the 'K' DO-loop (if desired; this is an option not actually used in this example).

```

KOUNT = 0
DO 280 K = 1, NDT

```

Compute the new radiometer body temperature distribution every tenth time increment.

```

      KOUNT = KOUNT + 1
      IF(KOUNT.EQ.10) GOTO 678
      GOTO 679
678  CONTINUE
      KOUNT = 0
219  CONTINUE
      DO 230 J = 1,NBODY
      CJ = CBODY(J)*VBODY(J)
      CJ = DTBODY/CJ
      SUM1 = 0.0D0
      SUM2 = 0.0D0
      SIGMA = SIG*AO/10000.0D0
      DO 220 I = 1,NBODY
      IF(I.EQ.J) GOTO 220
      FACTOR = EBODY(I)*SIG*ABODY(I)*DBODY(I,J)
      FACTOR = FACTOR*TGBODY(I)*TGBODY(I)*TGBODY(I)*TGBODY(I)
      FACTOR = FACTOR + GBODY(I,J)*TGBODY(I)
      SUM1 = SUM1 + FACTOR
      SUM2 = SUM2 + GBODY(I,J)
220  CONTINUE
      IF(J.GE.1.AND.J.LT.5)
&      QBODY(J) = SIGMA*SOURCE(5)*TSOURC**4
      IF(J.GE.5.AND.J.LT.9)
&      QBODY(J) = SIGMA*SOURCE(9)*TSOURC**4
      IF(J.GE.9.AND.J.LT.13)
&      QBODY(J) = SIGMA*SOURCE(13)*TSOURC**4
      IF(J.GE.17.AND.J.LT.21)
&      QBODY(J) = SIGMA*SOURCE(17)*TSOURC**4
      IF(J.GE.29.AND.J.LT.33)
&      QBODY(J) = SIGMA*SOURCE(21)*TSOURC**4
      IF(J.GE.33.AND.J.LT.37)
&      QBODY(J) = SIGMA*SOURCE(25)*TSOURC**4
      IF(J.GE.37.AND.J.LT.41)
&      QBODY(J) = SIGMA*(SOURCE(29)+SOURCE(33))*TSOURC**4
      IF(J.GE.41.AND.J.LT.45)
&      QBODY(J) = SIGMA*SOURCE(37)*TSOURC**4
      IF(J.GT.20.AND.J.LT.25) QBODY(J) = 3.46D0*TCLAMP
      IF(J.GT.24.AND.J.LT.29) QBODY(J) = 0.22D0*TCLAMP
      IF(J.GT.48.AND.J.LT.53) GOTO 636
      GOTO 637
636  CONTINUE
      ILL = J - 48
      QBODY(J) = QTOT(ILL)/10.0D0 + GLINK*T1BODY(J)
637  CONTINUE
      TOP = CJ*(SUM1 + QBODY(J)) + T1BODY(J)
      BOTTOM = (DBODY(J,J) - 1.0D0)*EBODY(J)*SIG*ABODY(J)
      BOTTOM = BOTTOM*TGBODY(J)*TGBODY(J)*TGBODY(J)
      IF(J.GT.20.AND.J.LT.25) SUM2 = SUM2 + 3.46D0
      IF(J.GT.24.AND.J.LT.29) SUM2 = SUM2 + 0.22D0
      IF(J.GT.48.AND.J.LT.53) SUM2 = SUM2 + GLINK
      BOTTOM = 1.0D0 + CJ*(SUM2 - BOTTOM)

```

```

TUBODY(J) = TOP/BOTTOM
IF(J.GT.76.AND.J.LT.81) TUBODY(J) = HST
230 CONTINUE
SUM = 0.0D0
AVE = 0.0D0
DO 240 I = 1,NBODY
BRMS = TGBODY(I) - TUBODY(I)
BRMS = BRMS*BRMS
SUM = SUM + BRMS
AVE = AVE + TGBODY(I)
240 CONTINUE
BRMS = SUM/RNBODY
BRMS = DSQRT(BRMS)
AVE = AVE/RNBODY
TEST = BRMS/AVE
IF(TEST.LE.0.000001) GOTO 241
DO 250 I = 1,NBODY
TGBODY(I) = TUBODY(I)
250 CONTINUE
GOTO 219
241 CONTINUE
DO 260 I = 1,NBODY
IF(I.LE.4) QTOT(I) = 0.0D0
T1BODY(I) = TUBODY(I)
260 CONTINUE
EETIME = ETIME + DT
WRITE(6,88) EETIME
DO 246 IS = 4, 104, 4
IX = IS - 3
WRITE(6,66) (T1BODY(I17), I17 = IX, IS)
246 CONTINUE
QBOD = 0.0D0
DO 661 I = 1,104
ALPHA = 0.96D0
IF(I.LT.13) ALPHA = 0.05D0
QBOD = QBOD
&      + ALPHA*ABODY(I)*SIG*DBODY(I,106)*T1BODY(I)**4
661 CONTINUE
QBOD = QBOD/4.0D0
679 CONTINUE

```

Compute the elapsed time.

```
ETIME = K*DT
```

The 'ITER' DO-loop loops through the temperature calculations until convergence is obtained.

```
DO 400 ITER = 1, ITRMAX
RMS = 0.0D0
TSUM = 0.0D0
RMSMAX = 0.0D0
```

The 'I' DO-loop steps through the nodes.

```
DO 18 I = 1, M2
```

Compute the (Level/Position) coordinates of Node I.

```
I1 = I - 1
I1 = I1/NTH
I2 = I - I1*NTH
I1 = I1 + 1
```

Compute the indices of the nodes surrounding Node I.

```
JJJ1 = I - 1
IF(I2.EQ.1) JJJ1 = JJJ1 + NTH
JJJ2 = I - NTH
JJJ3 = I + 1
IF(I2.EQ.NTH) JJJ3 = JJJ3 - NTH
JJJ4 = I + NTH
IF(JJJ2.LE.0) JJJ2 = I
IF(JJJ4.GT.M2) JJJ4 = I
```

Compute the heat flow rate radiated into each node during the calculation time interval DT.

```
Q1 = 0.0D0
DFSUM = 0.0D0
```

Loop through the J nodes which radiate to Node I.

```
DO 22 J = 1, M2
```

Compute (Level/Position) coordinates of Node J.

```
J1 = J - 1
J1 = J1/NTH
J2 = J - J1*NTH
J1 = J1 + 1
```

If Node J is in the first position on Level J1, i.e., if $J2 = 1$, the appropriate distribution factors have been read in; otherwise they must be found by a coordinate transformation.

```
IF(J2.EQ.1) GOTO 20
```

Coordinate transformation; rotate source Node J to first position on Level J1 and rotate receiver node an equal number of positions on Level I1.

```
JJ = J1
II = (I1 - 1)*NTH + I2 - J2 + 1
IF((I2 - J2).LT.0) II = II + NTH
GOTO 21
```

The distribution factor has been read in directly.

```
20 CONTINUE
JJ = J1
II = I
```

The distribution factor from Node J to Node I is the same as the distribution factor from Node JJ to Node II due to symmetry.

```
21 CONTINUE
```


Calculate Z(I,J).

```

IF(J.NE.I) GOTO 75
II1 = (I1 - 1)*NTH + 1
ZIJ = AA(I)*EC*SIG*TO(I)*TO(I)*TO(I)*DFC(II,II1)
Z(I,J) = -(EC + EB)*AA(I)*SIG*TO(I)*TO(I)*TO(I)
Z(I,J) = Z(I,J) + ZIJ - CC(I)/DT
Z(I,J) = Z(I,J) - G(JJJ1,I) - G(JJJ2,I)
Z(I,J) = Z(I,J) - G(JJJ3,I) - G(JJJ4,I)
GOTO 76
75 Z(I,J) = EC*SIG*TO(J)*TO(J)*TO(J)*AA(J)*DFC(JJ,II)
Z(I,J) = Z(I,J) + G(I,J)

```

There is no heat transfer by conduction from a node to to itself; thus if J = I, we skip the 'Q1' summation.

```

76 IF(J.EQ.I) GOTO 22
Q1 = Q1 + Z(I,J)*TN(J)
DFSUM = DFSUM + DFC(II,JJ)
22 CONTINUE
Q2 = 1.0D0 - DFSUM
Q2 = Q2*AA(I)*EC
Q2 = Q2*SIG*TSOURC*TSOURC*TSOURC*TSOURC

```

The heat input into the first ring of cavity nodes must include that from the F-O-V limiter; thus, if 'I' is less than or equal to 4, we must add the heat 'QBOD' from the F-O-V limiter.

```

IF(I.LE.4) Q2 = Q2 + QBOD

```

Calculate the quantity RHS(I).

```

RHS(I) = - AA(I)*EB*SIG*HST*HST*HST*HST
Q2 = Q2 + HEAT*AA(I)
RHS(I) = RHS(I) - Q2 - CC(I)*TO(I)/DT

```

Check to see if the matrix inversion method has been chosen; if so, proceed to next node.

```

IF(MAT.EQ.1) GOTO 888

```

Calculate TNEW for this node if the iterative solution has been chosen.

```

      IF(I.LE.4) GOTO 188
      GOTO 189
188  CONTINUE
      IHLP = I + 48
      TNEW = (Q1 - RHS(I) -
&          GLINK*T1BODY(IHLP))/(-Z(I,I) - GLINK)
      GOTO 190
189  CONTINUE
      TNEW = (Q1 - RHS(I))/(-Z(I,I))
190  CONTINUE
      TRMSMX = (TNEW - TN(I))*(TNEW - TN(I))
      RMS = RMS + TRMSMX
      IF(TRMSMX.GT.RMSMAX) RMSMAX = TRMSMX
      TSUM = TSUM + TNEW
      TN(I) = TNEW
      GOTO 18
1888 CONTINUE
      IF(I.LE.4) GOTO 517
      GOTO 518
1517 CONTINUE
      IHLP = I + 48
      RHS(I) = RHS(I) - GLINK*T1BODY(IHLP)
      Z(I,I) = Z(I,I) - GLINK
1518 CONTINUE
18  CONTINUE

```

If the matrix solution method has been chosen, call the matrix inversion Subroutine GAUSS.

```

      IF(MAT.EQ.0) GOTO 73
      CALL GAUSS(Z,RHS,M2,NARSZ,IAN5)
      IF(IER.EQ.131) STOP
      IF(IER.EQ.130) STOP
      DO 288 NP = 1, M2
      TN(NP) = RHS(NP)
288  CONTINUE
      GOTO 270

```

Determine if convergence has been obtained.

```

73 RMS = DSQRT(RMS)
   RMSMAX = DSQRT(RMSMAX)
   ERROR = RMS*100.0/TSUM
   ERRMAX = RMSMAX*100.0*FM2/TSUM
   IF((ERRMAX.LE.ELIMIT).AND.
     & (ITER.GT.ITRMIN)) GOTO 28
400 CONTINUE
410 CONTINUE
28 CONTINUE

```

Write the value of the elapsed time (s).

```

270 WRITE(6,8) ETIME

```

Write out current cavity temperature distribution at the end of each integration step.

```

DO 300 I5 = NTH, M2 , NTH
  I9 = I5 - NTH + 1
  WRITE(6,6) (TN(I3), I3 = I9, I5)
  WRITE(16,6) (TN(I3), I3 = I9, I5)
300 CONTINUE
DO 299 IJK = 1, M2
  TO(IJK) = TN(IJK)
  IJKL = IJK + 48
  IF(IJK.LE.4) LL = IJK
  IF(IJK.LE.4) QTOT(LL) = QTOT(LL)
  & + GLINK*(TO(IJK) - T1BODY(IJK))
299 CONTINUE
280 CONTINUE

```

Format statements.

```

1 FORMAT(5X, I5, 2D14.5)
2 FORMAT(5X, 3I5, D14.5)
3 FORMAT(5X, 2I5, 2D14.5)
4 FORMAT(5X, F10.4)
5 FORMAT(5X, 2I5, D14.5)
6 FORMAT(5X, 2(F8.3),2(F8.3),2(F8.3),
&      2(F8.3),2(F8.3))
7 FORMAT(5X, 2I5, 3D14.5)
8 FORMAT(10X, ///, 5X, 'ELAPSED TIME =',
&      F6.1, ' SECOND(S).', //)
9 FORMAT(5X/5X, 'CONVERGENCE NOT OBTAINED,',
&      ' ERROR =', E11.4, ' PERCENT'/)
66 FORMAT(45X, 2(F8.3),2(F8.3),2(F8.3),2(F8.3),2(F8.3))
88 FORMAT(44X, 'ELAPSED TIME =', F6.1, ' SECOND(S).', //)
81 FORMAT(4X, I3, 3X, F4.2, 1X, F4.2, 2X, F5.3, 5X, F5.3, 5X, F5.3)
82 FORMAT(4X, I3, 3X, F6.4)
83 FORMAT(5X, I2, 2X, I3, 2X, F5.2)
92 FORMAT(5X, 2I5, D14.5)
93 FORMAT(//5X, 'ITERATION MAXIMUM ERROR LIMIT ',
&      '= ', F10.8, ' PERCENT.', /, '1')
94 FORMAT('1', //15X, 'TRANSIENT',
&      'TEMPERATURE DISTRIBUTION OF A', /, 15X,
&      'TRW TYPE INVERTED CONE CAVITY RADIOMETER.',
&      //, 14X, 'DEVELOPED BY J. R. MAHAN AND ',
&      'L. D. ESKIN AT', /, 11X, 'VIRGINIA ',
&      'POLYTECHNIC INSTITUTE & STATE UNIVERSITY',
&      /, 17X, 'FOR THE NASA LANGLEY RESEARCH CENTER.')
95 FORMAT(10X, 'THE SOLUTION HAS SWITCHED ',
&      'TO THE ITERATION METHOD.')
96 FORMAT(5X, 'THE SYSTEM VARIABLES ARE AS ',
&      'FOLLOWS:' ,//10X, 'A = ', F5.2, ' MM',
&      /, 10X, 'B = ', F5.2, ' MM', /, 10X, 'C = ',
&      F5.2, ' MM', /, 10X, 'N1 = ', I2, /, 10X, 'N2 '
&      ', '= ', I2, /, 10X, 'NTH = ', I2, /, 10X, 'NSHOTN = '
&      ', I6, /, 10X, 'ABS = ', F3.1, /, 10X, 'REFR = '
&      ', '= ', I6, /, 10X, 'ABS = ', F3.1, /, 10X, 'REFR = '
&      ', F3.1, /, 10X, 'DT = ', F4.1, /, 10X, 'ELIMIT = '
&      ', E8.2, /10X, 'TEMPERATURE SOLUTION METHOD ',
&      ' (0=ITERATION,1=MATRIX) = ', I2)
97 FORMAT(F5.2, F5.2, F5.2, I2, I2, I2, I6,
&      F3.1, F3.1, F4.1, E8.2, 4I1)

```

```

98 FORMAT(////15X,'NUMBER OF CAVITY NODES = '
&      ,I3,/,15X,I2,' CAVITY NODE RINGS WITH ',I2,
&      ' NODES PER RING.',/,1X,
&      T27,'SYSTEM GEOMETRY',//1X,T41,'_____',/
&      ,1X,T27,'|<----2C---->| | |',1X,T27,'|',T40,
&      '| | |',/1X,T27,'|',T40,'| A |',/1X,T27,'|'
&      ,T40,'| | |',/1X,T27,'|',T40,'| V |',/'+',T42
&      ,'|',/1X,T28,'*',T39,'*',T44,'B',/1X,T29,'*',
&      T38,'*',T44,'|',/1X,T30,'*',T37,'*',T44,'|',/
&      ,1X,T31,'*',T36,'*',T44,'|',/1X,T32,'*',T35
&      ,'*',T44,'|',/1X,T33,'*',T34,'*',T44,'V',/
&      ,'+',T35,'_____',/,5X,'A = ',F5.2,
&      ' MM',5X,'B = ',F5.2,' MM',5X,'C = ',F5.2
&      , ' MM',/'1')
99 FORMAT(5X,/,5X,'THE AVERAGE RMS ERROR = '
&      ,F10.8,' PERCENT AFTER ',I2,' ITERATIONS .'
&      //,5X,'THE MAXIMUM RMS ERROR = ',F10.8,
&      ' PERCENT.',/,5X,'THE HEAT SINK ',
&      'TEMPERATURE IS ',F5.1,' K.',/)

```

End of 'TRW TEMP'.

```

STOP
END

```

4.3. Subroutine GAUSS

This is a standard matrix inversion subroutine which could be replaced at the discretion of the user if the argument list of the replacement routine was the same. We present this subroutine without further explanation.

4.3.1. Subroutine Listing

```

SUBROUTINE GAUSS(A,B,N,NARSZ, IANS)

IMPLICIT REAL*8(A-H,O-Z), INTEGER(I-N)
INTEGER IANS(NARSZ)
REAL*8 A(NARSZ,NARSZ),B(NARSZ)

COMMON /EMSGC/ IER

DO 10 I = 1,N
  IANS(I) = I
  AM = DABS(B(I))

DO 20 J = 1, N
  BM = DABS(A(I,J))
  IF ( BM .GT. AM ) AM = BM
20 CONTINUE

  IF(AM.EQ.0.0) GOTO 901

10 CONTINUE

  N1 = N + 1

DO 100 L = 1,N
  AMX = DABS(A(L,L))
  I = L
  J = L

DO 110 I1 = L,N

DO 120 J1 = L,N

  F1 = DABS(A(I1,J1))
  IF(AMX.GE.F1) GOTO 120
  AMX = F1
  I = I1
  J = J1

120 CONTINUE

110 CONTINUE

  IF(I.EQ.L) GOTO 140

DO 130 K = 1,N
  F1 = A(I,K)
  A(I,K) = A(L,K)
  A(L,K) = F1
130 CONTINUE

```

```
F1 = B(I)
B(I) = B(L)
B(L) = F1

140 CONTINUE

IF(J.EQ.L) GOTO 160

DO 150 K = 1,N
F1 = A(K,J)
A(K,J) = A(K,L)
A(K,L) = F1
150 CONTINUE

III = IANS(J)
IANS(J) = IANS(L)
IANS(L) = III

160 CONTINUE

F1 = A(L,L)
IF(F1.EQ.0.0) GOTO 902

DO 170 LP = 1,N
A(L,LP) = A(L,LP)/F1
170 CONTINUE

B(L) = B(L)/F1

IF(N.EQ.L) GOTO 200
N2 = L + 1

DO 180 M = N2,N

F1 = A(M,L)
IF(F1.EQ.0.0) GOTO 182

DO 185 LP = 1,N
A(M,LP) = A(M,LP)/F1 - A(L,LP)
185 CONTINUE

B(M) = B(M)/F1 - B(L)

182 CONTINUE

180 CONTINUE

100 CONTINUE

200 CONTINUE
```

```
F1 = 0.0
N3 = N-1

DO 210 LP = 1,N3
  L = N - LP
  F1 = 0.0
  N4 = L + 1

  DO 220 K = N4,N
    F1 = F1 + B(K)*A(L,K)
220 CONTINUE

  B(L) = B(L) - F1
210 CONTINUE

  DO 250 LP = 1,N
    A(1,LP) = B(LP)
250 CONTINUE

  DO 260 I = 1,N
    IA = IANS(I)
    B(IA) = A(1,I)
260 CONTINUE

  GOTO 999

901 IER = 130

  WRITE(6,666)
666 FORMAT(/,' GAUSS(F):  ZERO ROW INPUT')
  GOTO 999

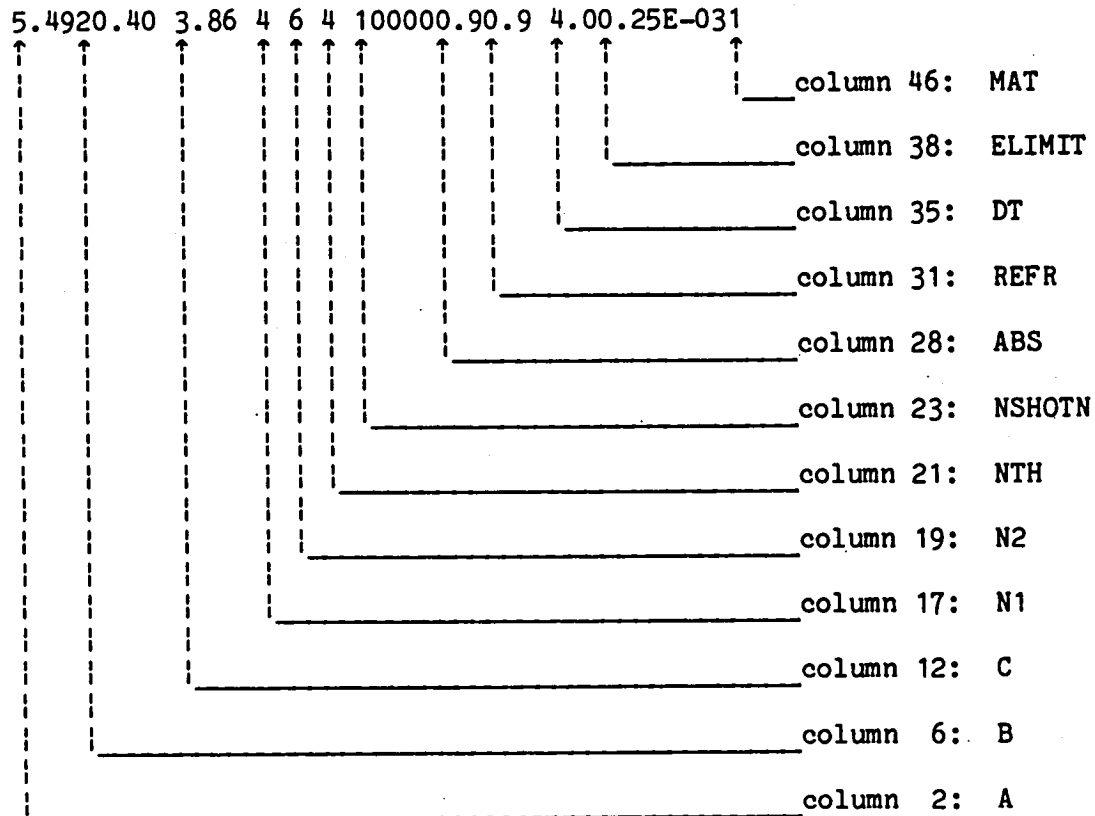
902 IER = 131

  WRITE(6,777)
777 FORMAT(/,' GAUSS(F):  SINGULAR MATRIX')
999 CONTINUE

RETURN
END
```


5. TRW DATA

The data read by 'TRW SHAPE', 'TRW NODES' and 'TRW TEMP' must be entered at the appropriate position at the end of each of the three programs. Note that not all of the programs use, or even read, all of the input data. However, the same data file serves all three.



<u>Variable</u>	<u>Type</u>	<u>Field</u>	<u>Description</u>
A	F5.2	1-5	Height of barrel (mm)
B	F5.2	6-10	Height of entire cavity (mm)
C	F5.2	11-15	Radius of barrel (mm)
N1	I2	16-17	Number of barrel axial divisions
N2	I2	18-19	Number of cone axial divisions
NTH	I2	20-21	Number of circumferential divisions
NSHOTN	I6	22-27	Number of volleys fired
ABS	F3.1	28-30	Absorptivity of cavity
REFR	F3.1	31-33	Reflectivity ratio
DT	F4.1	34-37	Cavity calculation time interval
ELIMIT	E8.2	38-45	Accuracy criterion using iteration
MAT	I1	46	= 1, matrix solution; = 0, iteration

6. BODY DATA

1	2.58	0.05	2.425	1.020	0.000
5	2.58	0.05	2.425	0.710	0.000
9	2.58	0.05	2.425	0.570	0.000
13	0.00	0.0	2.425	0.580	
17	3.05	0.96	2.425	1.900	0.000
21	8.24	0.0	2.425	9.080	
25	0.00	0.0	2.425	1.980	
29	3.90	0.96	2.425	1.330	0.000
33	1.56	0.96	2.425	1.200	0.000
37	1.20	0.96	2.425	0.120	0.000
41	0.24	0.96	2.425	0.260	0.000
45	0.46	0.0	2.425	0.530	
49	0.51	0.0	2.463	0.003	0.000
53	1.22	0.0	2.425	1.970	
57	0.56	0.0	2.463	0.003	
61	0.00	0.0	2.425	1.200	
65	0.40	0.0	3.430	0.240	
69	0.00	0.0	2.425	1.560	0.000
73	0.52	0.0	3.430	0.310	
77	0.00	0.0	3.430	1.540	0.000
81	0.51	0.0	3.430	0.310	
85	3.04	0.0	2.425	1.170	0.000
89	0.41	0.0	3.430	0.540	
93	4.87	0.0	2.425	1.340	
97	8.70	0.0	2.425	0.990	
101	1.07	0.0	2.463	0.016	

↑ column 41; QQBODY

↑ column 31; VVBODY

↑ column 21; CCBODY

↑ column 16; EEBOYD

↑ column 11; AABODY

↑ column 5; MBODY

<u>Variable</u>	<u>Type</u>	<u>Field</u>	<u>Description</u>
MBODY	I3	5-7	Body node number
AABODY	F4.2	11-14	Body node surface area
EEBODY	F4.2	16-19	Body node emissivity
CCBODY	F5.3	21-25	Body node thermal capacity
VVBODY	F5.3	31-35	Body node volume
QQBODY	F5.3	41-45	Body node heat input

These variables are described in detail in Section 4.1.

1	0.0592
5	0.0412
9	0.0320
13	0.0340
17	0.1384
21	0.6240
25	0.1240
29	0.1760
33	0.3440
37	0.0920
41	0.2000
45	0.4000
49	0.0192
53	0.8000
57	0.0208
61	0.4400
65	0.6800
69	0.5600
73	0.8800
77	1.0800
81	0.8800
85	3.9200
89	1.5600
93	6.7600
97	4.6400
101	1.0800

column 11; COND

column 5; MBODY

<u>Variable</u>	<u>Type</u>	<u>Field</u>	<u>Description</u>
MBODY	I3	5-7	Body node number
COND	F6.4	11-16	Conductance between body nodes on MBODY

1	5	8.95
2	6	8.95
3	7	8.95
4	8	8.95
5	9	6.59
6	10	6.59
7	11	6.59
8	12	6.59
9	13	5.32
10	14	5.32
11	15	5.32
12	16	5.32
13	17	11.23
14	18	11.23
15	19	11.23
16	20	11.23
13	21	2.08
14	22	2.08
15	23	2.08
16	24	2.08
17	21	7.20
18	22	7.20
19	23	7.20
20	24	7.20
17	29	3.91
18	30	3.91
19	31	3.91
20	32	3.91
21	25	4.33
22	26	4.33
23	27	4.33
24	28	4.33
29	33	2.63
30	34	2.63
31	35	2.63
32	36	2.63
33	37	0.66
34	38	0.66
35	39	0.66
36	40	0.66
33	41	0.02
34	42	0.02
35	43	0.02
36	44	0.02
33	45	0.04
34	46	0.04
35	47	0.04
36	48	0.04
33	53	0.05

34	54	0.05
35	55	0.05
36	56	0.05
37	41	0.00
38	42	0.00
39	43	0.00
40	44	0.00
41	45	7.11
42	46	7.11
43	47	7.11
44	48	7.11
45	53	4.03
46	54	4.03
47	55	4.03
48	56	4.03
49	57	0.03
50	58	0.03
51	59	0.03
52	60	0.03
53	61	7.77
54	62	7.77
55	63	7.77
56	64	7.77
57	65	0.06
58	66	0.06
59	67	0.06
60	68	0.06
61	65	1.99
62	66	1.99
63	67	1.99
64	68	1.99
61	69	8.44
62	70	8.44
63	71	8.44
64	72	8.44
65	73	3.17
66	74	3.17
67	75	3.17
68	76	3.17
69	73	2.59
70	74	2.59
71	75	2.59
72	76	2.59
69	77	8.92
70	78	8.92
71	79	8.92
72	80	8.92
73	81	2.82
74	82	2.82
75	83	2.82
76	84	2.82

77	81	4.51
78	82	4.51
79	83	4.51
80	84	4.51
77	85	13.46
78	86	13.46
79	87	13.46
80	88	13.46
81	89	2.06
82	90	2.06
83	91	2.06
84	92	2.06
85	89	9.17
86	90	9.17
87	91	9.17
88	92	9.17
85	93	0.72
86	94	0.72
87	95	0.72
88	96	0.72
89	101	0.03
90	102	0.03
91	103	0.03
92	104	0.03
93	97	2.47
94	98	2.47
95	99	2.47
96	100	2.47

column 16; GBODY(MIBODY,MJBODY)

column 10; MJBODY

column 6; MIBODY

<u>Variable</u>	<u>Type</u>	<u>Field</u>	<u>Description</u>
MIBODY	I2	6-7	Body node 'level' index
MJBODY	I3	10-12	Body node 'level' index
GBODY	F5.2	16-20	Conductance between MIBODY and MJBODY

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